GROUNDWATER TREATMENT SYSTEM QUARTERLY MONITORING REPORT SECOND QUARTER 2002

AMERICAN CHEMICAL SERVICE NPL SITE GRIFFITH, INDIANA

MWH File No. 2090601

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Prepared For:

American Chemical Service NPL Site RD/RA Executive Committee Griffith, Indiana

Prepared By:

MWH 27755 Diehl Road, Suite 300 Warrenville, Illinois 60555

February 2003





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Prepared For:

American Chemical Service NPL Site RD/RA Executive Committee Griffith, Indiana

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Approved by:	Peter Vagt, Ph.D., CPG Project Manager	February 7, 2009 Date

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- June 28, 2002 Off-Gas Sample (Round 4) Laboratory Results

NPL Site

1.0 INTRODUCTION

On behalf of the ACS RD/RA Executive Committee, MWH started up the groundwater treatment plant (GWTP) at the American Chemical Service NPL Site (ACS Site) in Griffith, Indiana on March 13, 1997. The GWTP was designed to treat groundwater from the Perimeter Groundwater Containment System (PGCS) and the Barrier Wall Extraction System (BWES). The original treatment consisted of a phase-separator for oil and free product removal, equalization tanks, a UV-oxidation unit for destruction of organic constituents, and an air stripper to remove methylene chloride and other organics. The treatment also included a chemical precipitation and clarification unit to remove metals, a sand filter to remove suspended solids, and activated carbon vessels for final polishing of the treated groundwater.

In 2001, an activated sludge treatment process was added to the process to reduce the volatile and semivolatile organic compounds (VOCs and SVOCs) in the collected groundwater. The activated sludge treatment process also reduces the amount of activated carbon required in the treatment process. An aerated equalization tank was also added to the GWTP in 2001 to remove VOCs from the collected groundwater, oxidize metals to increase metals removal efficiency in the chemical precipitation unit, and equalize groundwater flow through the GWTP. The activated sludge system and aeration tank have been fully integrated into the process along with the other upgrade components. Startup and optimization of the catalytic oxidizer/scrubber air treatment unit was also conducted during 2001.

MWH began eight initial rounds of off-gas sampling of the catalytic oxidizer/scrubber described in the PSVP during April 2002. The first four sampling rounds were conducted during the second quarter 2002.

The treated effluent from the treatment system is discharged to the nearby wetlands, west of the treatment system, in accordance with Agency approvals. This Groundwater Treatment System report summarizes effluent analytical data, catalytic oxidizer/scrubber off-gas analytical data, and water level gauging data collected from April 2002 through June 2002. This report details modifications and upgrades to the GWTP during the reporting period. This report also summarizes the sediment analytical data from the annual sediment sample collected during August 2001 at the wetlands discharge point of the GWTP.

2.0 COMPLIANCE MONITORING

2.1 INTRODUCTION

Effluent samples were collected from the treatment system to demonstrate compliance with the discharge limits (Table 2.1) established by Indiana Department of Environmental Management (IDEM) and United States Environmental Protection Agency (U.S. EPA). The approved Performance Standard Verification Plan (PSVP) requires quarterly effluent sampling for biological oxygen demand (BOD), total suspended solids (TSS), SVOCs, metals, and polychlorinated biphenyls (PCBs) in the system, and monthly effluent sampling for VOCs, as shown in the table below. To gather additional information, beyond the requirements of the PSVP, the effluent sampling was conducted on a monthly basis for all analytes. The samples will continue to be collected on a monthly basis until the treatment system is operating in a relatively steady state after completion and optimization of the groundwater treatment plant upgrades.

Sampling and analyses were performed in accordance with the approved Quality Assurance Project Plan (QAPP). Quality control measures were also instituted in accordance with the PSVP and QAPP. The following table and paragraphs present details on sampling and analyses and also summarize the analytical data for the treatment system effluent.

Sampling Frequency Schedule - Groundwater Treatment System

Analytes	Cumulative Time From Startup*	Frequency
Flowrate and pH	_	Continuous
BOD, TSS, SVOCs and Metals	181 days onward	Once per quarter
VOCs	31 days onward	Once per month
PCBs	181 days onward	Once per quarter
PCBs in Sediment (one location)		Once per year

^{*}Note: System was started up on March 13, 1997

2,2 EFFLUENT SAMPLING AND ANALYSES

Effluent samples were collected each month during the second quarter 2002. Samples were collected on the following dates for this reporting period:

- April 22, 2002
- May 9, 2002
- June 20, 2002

The above samples were collected directly from a sample tap on the effluent line of the treatment system. The samples were placed in contaminant-free containers, in accordance with the U.S. EPA Specifications and Guidance for Obtaining Contaminant-Free Sample Containers (U.S. EPA, 1992). Appropriate sample containers and preservatives, as specified in the QAPP, were used to collect and preserve the samples. Following sample collection, the sample containers were refrigerated at or below 4° C in coolers. Chain-of-

Custody forms were prepared to track the transfer of samples from the treatment system to the laboratories. In accordance with the approved QAPP, the effluent water samples were analyzed for the following parameters by the following analytical methods:

<u>Parameter</u>	Analytical Method
VOCs	SW-846 8260B
SVOCs	SW-846 8270C
Pentachlorophenol	SW-846 8270C and SIM
Pesticides/PCBs	EPA 608/SW-846 8081/8082
Metals (Excluding Mercury)	SW-846 6010
General Water Quality	EPA 160.2 and 405.1
Parameters (TSS and BOD-5)	
Mercury	SW-846 7470
рН	EPA 150.1

The sediment sample and associated quality control samples were analyzed for PCBs using analytical method SW-846-8082.

2.3 EFFLUENT ANALYTICAL RESULTS

GWTP Effluent Samples

The effluent monitoring data, summarized in Table 2.2, verifies that the system effluent was compliant with the discharge limits presented in Table 2.1. No exceedences were reported. The analytical data sheets for the compliance samples are provided in Appendix A.

Compuchem Laboratory of Cary, North Carolina performed the analysis of the samples. Laboratory Data Consultants (LDC) of Carlsbad, California performed third party data validation in accordance with the U.S. EPA National Functional Guidelines for Organic/Inorganic Data Review. Validation qualifiers are listed in Table 2.2 and are written in the margin of the analytical data sheets provided in Appendix A.

Sediment Sample

MWH conducted an investigation of the wetland areas north and west of the ACS site in May 1996 after earlier Remedial Investigations (RI) indicated the presence of PCBs. Locations for soil/sediment samples were selected by representatives of the U.S. EPA and MWH to more clearly delineate the extent and concentrations of PCBs in the wetland. Samples were collected from several locations across the wetlands. Results of this sampling indicated that low level residual PCBs, including Aroclor-1260, were present throughout the wetland. The wetland investigation is documented in the Phase I Technical Memorandum Wetland Investigation (MWH, July 1996) and the Phase II Technical Memorandum Wetland Investigation (MWH, February 1997). A summary of the investigation sampling results and a map of the sampling locations are included in Appendix B.

Since 1998, MWH has collected an annual sediment sample and associated quality control samples from the GWTP outfall in accordance with the PSVP to help determine if PCB accumulation is occurring at the GWTP discharge location. The annual sediment sample for 2002 was collected on June 5, 2002 from the GWTP outfall location shown on Figure 2.1. The sample was analyzed for PCBs by Compuchem and the data was validated by LDC.

Aroclor-1260 was detected in the sample (Compuchem reported an estimated concentration of 41 ug/kg) but not in the field duplicate sample (Compuchem reported non-detect with and 49 ug/kg detection limit). The sample was given a "J" flag by both Compuchem and LDC, indicating that the result was detected below the reporting limit and is an estimated concentration. This result is below the 1,000 ug/kg remediation objective used in the August 2001 PCB-Impacted Soil Wetland Excavation. No other PCBs were detected in either the sample or the field duplicate sample.

The estimated concentration of Aroclor-1260 in the sediment sample collected June 2002 was the first detection of Aroclor-1260 since annual sediment sampling began in 1998, though it was detected during the wetland investigation of May 1996. The field duplicate sample was collected as a split sample, however, and indicated a non-detect concentration of Aroclor-1260. The estimated concentration of Aroclor-1260 is also lower than the detection limit of previous annual samples where Aroclor-1260 has not been detected. This variability makes the trending of PCB concentrations at low levels difficult and inaccurate.

There have been no GWTP effluent exceedences of PCBs since the February 2000 sediment sample, demonstrating that there is little likelihood of PCBs accumulating in the wetlands due to GWTP discharge. MWH will continue to collect annual sediment samples from the GWTP outfall point, according to the PSVP, to monitor for any potential accumulation of PCBs.

A summary of the analytical data for the annual sediment samples, collected in December 1998, February 2000, August 2001, and June 2002 are summarized in Table 2.3. Analytical data for the June 2002 sample are included in Appendix C.

2.4 CATALYTIC OXIDIZER/SCRUBBER OFF-GAS SAMPLING AND RESULTS

Off-Gas Sampling

Influent and effluent off-gas samples were collected from the catalytic oxidizer/scrubber unit (ME-106) in the GWTP four times during the second quarter 2002. These samples consisted of the first four of eight rounds planned for the catalytic oxidizer/scrubber unit. Samples were collected on the following dates:

- April 26, 2002 (Round 1)
- May 22, 2002 (Round 2)
- June 21, 2002 (Round 3)
- June 28, 2002 (Round 4)

The samples were collected directly from a sample tap on the influent and effluent lines of the catalytic oxidizer/scrubber. One influent sample (labeled IN1) and one effluent sample (EF1) were collected. A duplicate influent sample (IN2) was also collected. The samples were collected in accordance with the QAPP and laboratory guidelines. The VOC sample was collected using a summa canister and the SVOC sample was collected in sorbent tubes.

Following sample collection, the SVOC sample containers were refrigerated at or below 4°C in coolers. The VOC samples to be analyzed by method TO-14 do not require refrigeration. Chain-of-Custody forms were prepared to track the transfer of samples from the treatment system to the laboratories for extraction and analysis. In accordance with the approved QAPP, the off-gas samples were analyzed for the following parameters by the following analytical methods:

<u>Parameter</u>	Analytical Method
VOCs	TO-14
SVOCs	TO-13

Sampling Results

The influent and effluent off-gas data, summarized in Tables 2.4 and 2.5, verifies that the off-gas from the catalytic oxidizer was less than the IDEM discharge limit of 3 pounds per hour VOC discharge. For example, the VOC discharge reported from the April 26, 2002 sample was 0.03 pounds per hour, approximately one percent of the discharge limit. The analytical data sheets for the compliance samples are provided in Appendix D.

Air Toxics Laboratories of Folsom, California analyzed the samples. The analytical results are summarized in Tables 2.4 and 2.5. MWH performed data validation in accordance with the QAPP and the National Functional Guidelines for Organic/Inorganic Data Review. Validation qualifiers are listed in Tables 2.4 and 2.5 and are written in the margin of the analytical data sheets provided in Appendix D.

3.0 TREATMENT SYSTEM PROCESS MODIFICATIONS

The thermal oxidizer/scrubber unit, which is housed inside and next to the GWTP, began operating in April 2002 as part of the Off-Site Area In-Situ Soil Vapor Extraction (ISVE) system. The thermal oxidizer/scrubber unit is not part of the groundwater treatment process; however, it does interact with the GWTP. The scrubber utilizes the treated effluent water from the GWTP to maintain a sufficient water level. By doing so, the amount of city water usage is significantly decreased. This was accomplished by installing a new pump and piping from process tank T-1 to the scrubber. Operation of the pump is interlocked with the catalytic oxidizer programmable logic controller (PLC) and the water level sensors in tank T-1. The pump is enabled based on scrubber demand and an allowable water volume in tank T-1.

The scrubber periodically discharges water to the influent of GWTP for treatment to control the conductivity of the scrubber water. There are two factors that contribute to an increase in the conductivity, the continuous addition of sodium hydroxide and its subsequent reaction with the chlorides in the vapor stream to form sodium chloride. Once the conductivity, which is measured continuously, rises above a setpoint, the scrubber "blows down" a portion of the scrubber water. The "blowdown" is pumped to the gravity phase separator, ME-101, for treatment.

Performance and compliance testing of the thermal oxidizer and scrubber system are performed in conjunction with performance monitoring of the ISVE system and is, therefore, not included in this report.

4.0 PGCS AND BWES GAUGING ACTIVITIES

The PGCS trench groundwater extraction wells were operated in "auto" mode continuously throughout the second quarter 2002. In "auto" mode, each of the PGCS extraction wells is set to turn on or off automatically based on water levels within the Aeration Equalization Tank (T-102) and the individual extraction wells. This mode is used to control the flowrate through the treatment system. The GWTP also received influent from the BWES during the second quarter 2002.

MWH continued to regularly monitor water levels inside the barrier wall. Figure 4.1 shows the water levels as measured on June 28, 2002. Piezometers P29, P31, P32, P36, P49, P106, and P108 in the On-Site Area and P96, P110, P112, P113, P114, P116, and P118 in the Off-Site Area were measured regularly throughout the quarter. The water levels from these piezometers are listed in Table 4.1 and are depicted graphically on Figures 4.2 and 4.3. The target water levels in each area are shown on these figures for reference.

In late 2001, MWH revised the long-term groundwater monitoring program. The revisions included changing the frequency of collecting groundwater levels and analyzing groundwater samples from quarterly basis to a semi-annual basis. In the past, groundwater levels were collected from piezometers outside the barrier wall for the Groundwater Treatment System Quarterly Monitoring Report at the same time as groundwater levels for the groundwater monitoring program. Because of the change in the groundwater monitoring program, water levels from piezometers outside the barrier wall were not collected for the second quarter 2002. Thus, several water level measurements specific to the PSVP, including several piezometers in and around the PGCS, were not collected. This omission of data points specific to the PSVP was not noticed until the third quarter 2002. MWH then updated the LTGMP to include measurement of all PSVP water level locations on a quarterly schedule.

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Table 2.1 Groundwater Treatment System Effluent Discharge Limits American Chemical Service NPL Site Griffith, Indiana

Groundwater Quality Parameter	Effluent Standard (Limit)			
General Water Quality Parameters				
РН	6 - 9 S.U.			
BOD-5	30 mg/L			
TSS	30 mg/L			
Inorganics				
Arsenic	50 μg/L			
Beryllium	NE			
Cadmium	4.1 μg/L			
Manganese	NE			
Mercury	$0.02 \mu g/L (w/DL = 0.64)$			
Selenium	8.2 μg/L			
Thallium	NE			
Zinc	411 μg/L			
Volatile Organics				
Acetone	6.800 µg/L			
Benzene	5 μg/L			
2-Butanone	210 μg/L			
Chloromethane	NE			
1,4 – Dichlorobenzene	NE NE			
1,1 - Dichloroethane	NE			
1,2 – Dichloroethene – cis	70 μg/L			
Ethylbenzene	34 μg/L			
Methylene chloride	5 μg/L			
Tetrachloroethene	5 μg/L			
Trichloroethene	5 μg/L			
Vinyl chloride	2 μg/L			
4 – Methyl - 2 – pentanone	15 μg/L			
Semi-Volatile Organics				
bis(2 - Chloroethyl) ether	9.6 μg/L			
bis(2 - Ethylhexyl) phthalate	6 μg/L			
Isophorone	50 μg/L			
4 – Methylphenol	34 μg/L			
Pentachlorophenol	1 μg/L			
PCBs				
PCBs	$0.00056 \mu\text{g/L} (\text{w/DL} = 0.1 \text{ to } 0.9)$			

Notes:

NE = No effluent limit established.

DL = Detection limit

Table 2.2

Summary of Effluent Analytical Results - Second Quarter 2002 Groundwater Treatment System American Chemical Service NPL Site Griffith, Indiana

Event	Month 59	Month 60	Month 61	Effluent Limits	Lab Reporting
Date	4/22/02	5/9/02	6/20/02	Ettibent Chinis	Limits
pH	7.96	7.83	7.13 /J	6-9	none
TSS	ND	ND	1.2	30	10
BOD	ND	ND	24	30	2
Arsenic	ND	ND	ND	50	3.4
Beryllium	ND	0 21 B/UB	ND	NE	0.2
Cadmium	ND	ND	ND	4.1	0.3
Manganese	14.4	8.9 B/	24.5	NE	10
Mercury	ND	ND	ND	0.02 (w/DL = 0.64)	0.64
Selenium	ND	ND	ND	8.2	4.3
Thallium	ND	ND	ND	NE	5.7
Zinc	7.1 B/UB	ND	3.9 B/	411	1.2
Benzene	ND	ND	ND	5	0.5
Acetone	1 JB/	ND /UJ	4 B/UBJ	6,800	3
2-Butanone	ND	ND	ND /UJ	210	3
Chloromethane	0.1 J/3 UBJ	ND	0.2 3/J	NE	0.5
1,4-Dichlorobenzene	ND	ND	ND	NE	0.5
1,1-Dichloroethane	ND	ND	ND	NE	0.5
cis-1,2-Dichloroethene	ND	ND	ND	70	0.5
Ethylbenzene	ND	ND	ND	34	0.5
Methylene chloride	0.1 J/	ND	I B/UBJ	5	0.6
Tetrachloroethene	0.05 JB/0.5 UB	ND	ND	5	0.5
Trichloroethene	ND	ND	ND	5	0.5
Vinyl chloride	ND	ND_	ND	2	0.5
4-Methyl-2-pentanone	ND	ND /UJ	ND	15	3
bis (2-Chloroethyl) ether	ND	ND	ND	9.6	9.6
bis(2-Ethylhexyl) - phthalate	0.91 JB/6 UB	ND	4 JB/UB	6	6
4 - Methylphenol	ND	_ND	ND	34	10
lsophoron e	ND	ND /UJ	ND	50	10
Pentachlorophenol	ND	ND	ND	1	1
PCB/Aroclor-1016	ND	ND	ND	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1221	ND	ND	ND	0.00056 (w/DL = 0.1 to 0.9)	0.92*
PCB/Aroclor-1232	ND	ND	ND	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1242	ND	ND	ND /UJ	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1248	ND	ND	ND /UJ	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1254	ND	ND	ND /UJ	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1260	ND	ND	ND /UJ	0.00056 (w/DL = 0.1 to 0.9)	0.5

Notes

Data has been validated in accordance with the Project QAPP (November 2001) and the U.S. EPA

National Functional Guidelines for Organic Data Review

Shaded cells indicate discharge exceedances

 $pH\ data$ is expressed in $S\ U$

TSS and BOD5 data is expressed in mg/L

Metals, VOC, SVOC and PCB data is expressed in ug/L

ND = Not detected

NE = No effluent limit established

NA = Sample not analyzed for this compound

* = Approved SW 846 method is incapable of achieving effluent limit

Suffix Definitions

- _/ = Data qualifier added by laboratory
- /_ = Data qualifier added by data validator
- B = Compound is also detected in the blank
- E = Compound exceeds the upper level of calibration range of instrument
- J = Result is detected below the reporting limit and is an estimated concentration
- JB = Analyte is detected in the compliance sample below the reporting limit and is an estimated concentration and the compound is also detected in the method blank resulting in a potential high bias
- $\mathbf{U} = \mathbf{A}\mathbf{n}\mathbf{a}\mathbf{l}\mathbf{y}\mathbf{t}\mathbf{e}$ is not detected at or above the indicated concentration
- $UB\cong Analyte$ is not detected at or above the indicated concentration due to blank contamination
- UJ = Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.

Table 2.3 Summary of Sediment Analytical Results Groundwater Treatment System American Chemical Service NPL Site Griffith, Indiana

DCD C	Results (ug/kg)										
PCB Compound	12/4/98	2/3/00	2/3/00 DUP	8/21/01	8/21/01 DUP	6/5/02	6/5/02 DUP				
Aroclor-1016	ND (33)	ND (59)	ND (79)	ND (62) /UJ	ND (71)	ND (52) /UJ	ND (49)				
Aroclor-1221	ND (33)	ND (77)	ND (100)	ND (82) /UJ	ND (92)	ND (67) /UJ	ND (64)				
Aroclor-1232	ND (33)	ND (59)	ND (79)	ND (62) /UJ	ND (71)	ND (52) /UJ	ND (49)				
Aroclor-1242	ND (33)	ND (41)	ND (55)	ND (43) /UJ	ND (49) /UI	ND (36) /UJ	ND (34)				
Aroclor-1248	ND (33)	ND (41)	ND (55)	ND (43) /UJ	ND (49) /UI	ND (36) /UJ	ND (34)				
Aroclor-1254	ND (33)	22 J/	15 J/	73 P/J	39 JP/J	ND (36) /UJ	ND (34)				
Aroclor-1260	ND (33)	ND (59)	ND (79)	ND (62) /UJ	ND (71) /UJ	41 3/3	ND (49)				
Total PCBs	ND	22	15	7.3	39	41	ND				

Notes

- 1 ND (_) = Compound was not detected. The detection limit is included in parentheses
- 2 December 4, 1998 sample was analyzed by Quanterra All other samples were analyzed by Compuchem
- 3. DUP = Duplicate sample

Suffix Definitions

- _/ = Data qualifier added by laboratory
- I_{\perp} = Data qualifier added by data validator
- B = Compound is also detected in the blank
- J = Result is detected below the reporting limit and is an estimated concentration
- P = The Relative Percent Difference (RPD) between the two GC column values is greater than 25%. The higher value has been reported.
- JB = Analyte is detected in the sample below the reporting limit and is an estimated concentration. The compound is also detected in the method blank resulting in a potential high bias
- UB = Analyte is not detected at or above the indicated concentration due to blank contamination
- UI = Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- JP = Result is detected below the reporting limit and is an estimated concentration.

 Also, the Relative Percent Difference (RPD) between the two GC column values is greater than 25%. The higher value has been reported.

American Chemical Service, Griffith, Indiana

	T	Round 1 - Sampled 4/26/02							
	ľ		Analytical Dat	Destruction Efficiency					
Compounds	Units	Influent IN1	Influent IN2	Effluent EF1	Low	High	Average		
Method TO-14	Г								
Chloromethane	ppbv	ND	ND	240	NC	NC	NC		
Vinyl Chlonde	ppbv	2,900	7,300	350	87.93%	95.21%	91.57%		
Bromomethane	ppbv	ND	ND	ND	NC	NC	NC		
Chioroethane	ppbv	2,200	6,600	170	92.27%	97.42%	94.85%		
1,1-Dichloroethene	ppbv	24]/J	410	24	NC	94.15%	94 15%		
Methylene Chloride	ppbv	3,100	88,000	440	85.81%	99.50%	92.65%		
1,1-Dichloroethane	ppbv	1,000	20,000	74	92.60%	99.63%	96 12%		
cis-1,2-Dichloroethene	ppbv	10,000	9,700	880	90.93%	91.20%	91 06%		
Chlorokform	ppbv	ND	2,000	ND	100 00%	100.00%	100.00%		
1,1.1-Trichloroethane	ppbv	270	53,000	13	95.19%	99.98%	97 58%		
Carbon Tetrachloride	ppbv	ND	ND	ND	NC	NC	NC		
Benzenc	ppov	14,000	33,000	1,500	89.29%	95.45%	92.37%		
1,2-Dichloroethane	ppbv	310	1,000	32	89.68%	96.80%	93.24%		
Frichloroethene	ppbv	65	18,000	7.1	89.08%	99.96%	94.52%		
1,2-Dichloropropane	ppbv	110	270 J/J	7.4	93.27%	NC	93.27%		
cis-1,3-Dichloropropene	ppbv	ND	ND	ND	NC	NC	NC		
Toluene	ppbv	10,000	100,000	750	92.50%	99.25%	95.88%		
rans-1,3-Dichloropropene	ppbv	ND	ND	ND	NC	NC	NC		
1.1.2-Trichloroethane	ppbv	20 J/J	ND	2.0 J/J	NC	NC	NC		
Tetrachloroethene	ppbv	23 J/J	6,800	5.3 J/J	NC	NC	NC		
Chlorobenzene	ppbv	920	ND	110	NC	88.04%	88.04%		
Ethylbenzene	ppbv	1,600	7,100	100	93.75%	98.59%	96.17%		
n,p-Xylene	ppbv	7,700	26,000	480	93.77%	98.15%	95.96%		
o-Xylene	ppbv	2,500	7,300	160	93.60%	97.81%	95.70%		
Styrene	ppbv	ND	ND	19	NC	NC	NC		
1,1,2,2-Tetrachloroethane	ppbv	ND	ND	ND	NC	NC	NC		
Acetone	ppbv	610	16,000	88	85.57%	99.45%	92.51%		
Carbon Disulfide	ppbv	ND	ND	ND	NC	NC	NC		
rans-1.2-Dichloroethene	ppbv	ND	ND	54	NC	NC	NC		
2-Butanone (MEK)	ppbv	340	15,000	27	92.06%	99.82%	95.94%		
Bromodichloromethane	ppbv	ND	ND	ND	NC	NC	NC		
1-Methyl-2-pentanone	ppbv	410	3,800	21 J/J	NC	NC	NC		
2-Hexanone	ppbv	ND	ND	ND	NC	NC	NC		
Dibromochloromethane	ppbv	ND	ND	ND	NC	NC	NC		
Bromoform	ppbv	ND	ND	ND	NC	NC	NC		
	L.,		404.046		00.400	00.00	04.505		
Total	ppbv	58,035	421,010	5,526	90.48%	98.69%	94.58%		
Total	lb/hr	0.324	2.527	0.030	NC	NC	NC_		
DEM Discharge Requirement	lb/hr	L		3					

Notes:

Qualifiers: J - Result is estimated

/ - Laboratory data qualifier

/_ - Data validation qualifier

NC - Not calculated

ND - Non-detect

ppby - parts per billion volume

Destruction efficiency is not calculated where influent and/or effluent values are estimated

Total VOCs in lb/hr calculation based on air flow rate of V = 400 acfm

American Chemical Service, Griffith, Indiana

	Ī	Round 2 - Sampled 5/22/02						
i	ſ	/	nalytical Dat	Destruction Efficiency				
Compounds	Units	Influent IN1	Influent IN2	Effluent EF1	Low	High	Average	
Method TO-14						<u> </u>		
Chloromethane	ppbv	ND	ND	17	NC	NC	NC	
Vinyl Chloride	ppbv	540	410	57	86.10%	89.44%	87.77%	
Bromomethane	ppbv	ND	ND	ND	NC	NC	NC	
Chloroethane	ppbv	280	210	18	91.43%	93.57%	92.50%	
1,1-Dichloroethene	ppbv	4.9]/]	3.7 J/J	6.1	NC	NC	NC	
Methylene Chloride	ppbv	420	360	49	86.39%	88.33%	87.36%	
I,I-Dichloroethane	ppbv	210	170	14	91.76%	93.33%	92.55%	
cis-1,2-Dichloroethene	ppbv	2,300	1,900	180	90 53%	92.17%	91 35%	
Chlororform	ppbv	3.8 J/J	3.1 1/3	0.54 J/J	NC	NC	NC	
1.1.1-Trichloroethane	ppbv	54	40	3.1	92.25%	94.26%	93 25%	
Carbon Tetrachloride	ppbv	ND	1.9 1/1	ND	NC	NC	NC	
Benzene	ppbv	2,200	1,800	240	86.67%	89.09%	87.88%	
1,2-Dichloroethane	ppbv	53	46	5.4	88.26%	89.81%	89 04%	
Trichloroethene	ppbv	6.3 J/J	5.4 3/3	0.8	NC	NC	NC	
1,2-Dichloropropane	ppbv	22	20	1.4	93.00%	93.64%	93.32%	
cis-1,3-Dichloropropene	ppbv	ND	ND	ND	NC	NC	NC	
Toluene	ppbv	1,700	1,400	120	91.43%	92.94%	92.18%	
trans-1,3-Dichloropropene	pphy	ND	ND	ND	NC	NC	NC	
1.1.2-Trichloroethane	ppbv	7.2 J/J	6.0 J/J	0.61 J/J	NC	NC	NC	
Tetrachloroethene	ppbv	ND	ND	0.30 J/J	NC	NC	NC	
Chlorobenzene	ppbv	200	170	23	86.47%	88.50%	87.49%	
Ethylbenzene	ppbv	410	350	20	94.29%	95.12%	94.70%	
m.p-Xylene	ppbv	1,700	1,400	74	94.71%	95.65%	95.18%	
o-Xylene	ppbv	580	520	26	95.00%	95.52%	95.26%	
Styrene	ppbv	ND	ND	3.8	NC	NC	NC	
1.1.2.2-Tetrachloroethane	ppbv	3.5 J/J	2.7 3/3	0.33 J/J	NC	NC	NC	
Acetone	ppbv	1,100	1,100	92	91.64%	91.64%	91.64%	
Carbon Disulfide	ppbv	ND	ND	0.31 J/J	NC	NC	NC	
trans-1,2-Dichloroethene	ppbv	ND	ND	8.5	NC	NC	NC	
2-Butanone (MEK)	ppbv	630	630	34	94.60%	94.60%	94.60%	
Bromodichloromethane	ppbv	ND	ND	ND	NC	NC	NC	
4-Methyl-2-pentanone	ppbv	440	440	15	96.59%	96.59%	96.59%	
2-Hexanone	ppbv	13 J/J	12 J/J	0.46 J/J	NC	NC	NC	
Dibromochloromethane	ppbv	ND	ND	ND	NC	NC	NC	
Bromoform	ppbv	ND/UJ	ND /UJ	ND /UJ	NC _	NC	NC	
Total	ppbv	12,839	10,966	1.008	92.15%	90.81%	91.48%	
Total	lb/hr	0.069	0.059	0.005	NC	NC NC	NC	
IDEM Discharge Requirement	lb/hr	0.007	0.037	3	110	ITC	110	

Notes:

Qualifiers:

J - Result is estimated

/ - Laboratory data qualifier / - Data validation qualifier

NC - Not calculated

ND - Non-detect

ppbv - parts per billion volume

Destruction efficiency is not calculated where influent and/or effluent values are estimated.

Total VOCs in lb/hr calculation based on air flow rate of V = 390 acfm

Table 2.4

Summary of Catalytic Oxidizer Off-Gas Analytical Results for VOCs (Method TO-14) Second Quarter 2002

American Chemical Service, Griffith, Indiana

	Round 3 - Sampled 6/21/02							
}	1	,	Analytical Dat	Destruction Efficiency				
Compounds	Units	Influent INI	Influent IN2	Effluent EF1	Low	High	Average	
Method TO-14								
Chloromethane	ppbv	ND	ND	170	NC	NC	NC	
Vinyl Chloride	ppbv	2,600	2,600	290	88.85%	88 85%	88.85%	
Bromomethane	ppbv	ND	ND	ND	NC	NC	NC	
Chloroethane	ppbv	1,700	1,700	130	92.35%	92.35%	92.35%	
1,1-Dichloroethene	ppby	28 J/J	23 J/J	26 J/J	NC	NC	NC	
Methylene Chloride	ppbv	1,700	1,600	210	86.88%	87.65%	87.26%	
1,1-Dichloroethane	ppbv	890	860	54	93.72%	93.93%	93.83%	
cis-1,2-Dichloroethene	ppbv	9,400	9,200	750	91 85%	92 02%	91.93%	
Chlororform	ppbv	15 3/J	ND	ND	NC	NC	NC	
1,1,1-Trichloroethane	ppbv	300	290	12 J/J	NC	NC	NC	
Carbon Tetrachloride	ppbv	ND	ND	ND	NC	NC	NC	
Benzene	ppbv	15,000	15,000	1,400	90 67%	90.67%	90.67%	
1,2-Dichloroethane	ppbv	ND	ND	ND	NC	NC	NC.	
Trichloroethene	ppbv	63 J/J	60 J/J	9.0 J/J	NC	NC	NC	
1,2-Dichloropropane	ppbv	78	78	7.7 J/J	NC	NC	NC	
cis-1,3-Dichloropropene	ppbv	ND	ND	ND	NC	NC	NC	
Toluene	ppbv	8,500	8,400	570	93.21%	93.29%	93.25%	
trans-1.3-Dichloropropene	ppbv	ND	ND	ND	NC	NC	NC	
1.1.2-Trichloroethane	ppby	ND	ND	ND	NC	NC	NC	
Tetrachloroethene	ppbv	19 3/3	17 J/J	5 3 1/1	NC	NC	NC	
Chlorobenzene	ppbv	740	750	89	88.13%	87.97%	88.05%	
Ethylbenzene	ppbv	1,700	1,700	84	95 06%	95.06%	95.06%	
m,p-Xylene	ppbv	8,900	8,700	360	95.86%	95.96%	95.91%	
o-Xylene	ppbv	2,800	2,700	120	95.56%	95.71%	95.63%	
Styrene	ppbv	ND	ND	22 1/1	NC	NC	NC	
1,1,2,2-Tetrachloroethane	ppbv	ND	ND	ND	NC	NC	NC	
Acetone	ppbv	1,200	1,100	200	81.82%	83.33%	82.58%	
Carbon Disulfide	ppbv	ND	ND	ND	NC	NC	NC	
trans-1,2-Dichloroethene	ppbv	60 J/J	ND	86 J/J	NC	NC	NC	
2-Butanone (MEK)	ppbv	610	580	8.300	NC	NC	NC	
Bromodichloromethane	ppbv	ND	ND	ND	NC	NC	NC	
4-Methyl-2-pentanone	ppbv	490	420	22 J/J	NC	NC	NC	
2-Hexanone	ppbv	ND	ND	ND	NC	NC	NC	
Dibromochloromethane	ppbv	ND	ND	ND	NC	NC	NC	
Bromoform	ppbv	ND	ND	ND	NC	NC	NC	
Total	ppbv	56,608	55,678	12,727	77.14%	77.52%	77.33%	
Total	lb/hr	0.254	0.249	0.049	NC	NC	NC	
IDEM Discharge Requirement	lb/hr			3				

Qualifiers: J - Result is estimated

Notes:
_/ - Laboratory data qualifier
/_ - Data validation qualifier

NC - Not calculated

ND - Non-detect

ppby - parts per billion volume

Destruction efficiency is not calculated where influent and/or effluent values are estimated

Total VOCs in lb/hr calculation based on air flow rate of V = 320 acfm

American Chemical Service, Griffith, Indiana

	1	Round 4 - Sampled 6/28/02						
Í	i		Analytical Dat	a	Destr	uction Eff	iciency	
Compounds	Units	Influent IN1	Influent IN2	Effluent EF1	Low	High	Average	
Method TO-14	1							
Chloromethane	ppbv	ND	ND	180	NC	NC	NC	
Vinyl Chlonde	ppbv	3,900	4,600	440	88.72%	90.43%	89 58%	
Bromomethane	ppbv	ND	ND	ND	NC	NC	NC	
Chloroethane	ppbv	2,000	2,300	170	91.50%	92 61%	92 05%	
1,1-Dichloroethene	ppbv	25 J/J	29 J/J	32	NC	NC	NC	
Methylene Chloride	ppbv	860	1,200	130	84.88%	89.17%	87.03%	
1,1-Dichloroethane	ppbv	780	970	58	92.56%	94.02%	93.29%	
cis-1,2-Dichloroethene	ppbv	7,700	10,000	810	89 48%	91.90%	90.69%	
Chlororform	ppbv	12 J/J	13 J/J	1 2 J/J	NC	NC	NC	
1.1.1-Trichioroethane	ppbv	340	410	15	95.59%	96 34%	95 96%	
Carbon Tetrachloride	ppbv	ND	ND	ND_	NC	NC	NC	
Benzene	ppbv	14,000	17,000	(1,700)	87.86%	90.00%	88.93%	
1,2-Dichloroethane	ppbv	ND	ND	ND	NC	NC	NC	
Trichloroethene	ppbv	68	89	8.4	87.65%	90 56%	89.10%	
1.2-Dichloropropane	ppbv	60 J/J	87	4.5 J/J	NC	NC	NC	
cis-1.3-Dichloropropene	ppbv	ND	ND	ND	NC	NC	NC	
Toluene	ppbv	7,600	12,000	640	91.58%	94.67%	93 12%	
trans-1,3-Dichloropropene	ppbv	ND	ND	ND	NC	NC	NC	
1.1.2-Trichloroethane	ppbv	ND	ND	ND	NC	NC	NC	
Tetrachloroethene	ppbv	31 J/J	31 J/J	6.9 J/J	NC	NC	NC	
Chiorobenzene	ppbv	570	820	80	85.96%	90.24%	88.10%	
Ethylbenzene	ppbv	1,600	2,300	93	94.19%	95.96%	95.07%	
m.p-Xylene	ppbv	8,400	12,000	460	94.52%	96.17%	95.35%	
o-Xylene	ppbv	2,200	3,300	130	94.09%	96.06%	95.08%	
Styrene	ppbv	63 J/J	ND	22	NC	NC	NC	
1,1,2,2-Tetrachloroethane	ppbv	ND	ND	ND	NC	NC	NC	
Acetone	ppbv	350	400	48	86.29%	88.00%	87.14%	
Carbon Disulfide	ppbv	ND	ND	7.2 J/J	NC	NC	NC	
trans-1,2-Dichloroethene	ppbv	ND	ND	54	NC	NC	NC	
2-Butanone (MEK)	ppbv	200 J/J	240 J/J	32	NC	NC	NC	
Bromodichloromethane	ppbv	ND	ND	ND	NC	NC	NC	
4-Methyl-2-pentanone	ppbv	240 J/J	330	10 J/J	NC	NC	NC	
2-Hexanone	ppbv	ND	ND	ND	NC	NC	NC	
Dibromochloromethane	ppbv	ND	ND	ND	NC	NC	NC	
Bromoform	ppbv	ND	ND	ND	NC_	NC_	NC	
Total	ppbv	50,368	67,806	5,102	89.87%	92.48%	91.17%	
Total	lb/hr	0.240	0.325	0.023	NC	NC	NC	
IDEM Discharge Requirement	lb/hr			3				

high

Notes:

Qualifiers:

J - Result is estimated

_/ - Laboratory data qualifier

/_ - Data validation qualifier

NC - Not calculated

ND - Non-detect

ppby - parts per billion volume

Destruction efficiency is not calculated where influent and/or effluent values are estimated.

Total VOCs in $\frac{1}{V}$ acculation based on air flow rate of V = 340 acfm

American Chemical Service, Griffith, Indiana

				ound 1 - Sample		-4	Efficience		
	144.4		Analytical Data			struction Ef			
Compounds	Units	Influent IN1	Influent IN2	Effluent EF1	Low (%)	High (%)	Average (%		
Method TO-13		1/10	215	- NB	NO	NC	100		
Phenol	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC		
bis(2-Chloroethyl)ether 2-Chlorophenol	μg	ND ND	ND	ND	NC NC	NC NC	NC NC		
1.3-Dichlorobenzene	μg	0.88 J/J	ND	ND	NC NC	NC NC	NC NC		
1,4-Dichlorobenzene	μg	12	ND	1.0	NC NC	91.67%	91.67%		
1,2-Dichlorobenzene	μg	21	3.9	1.6	58.97%	92.38%	75.68%		
2-Methylphenol (o-Cresol)	μв	ND	ND	ND	NC	92.38% NC	73.06% NC		
N-Nitroso-di-n-propylamine	μg	ND	ND	ND ND	NC NC	NC NC	NC NC		
4-Methylphenol	μg	ND	ND	ND	NC	NC	NC		
Hexachloroethane	μg μg	ND ND	ND	ND	NC NC	NC	NC		
Nitrobenzene	μв	ND	ND	ND	NC NC	NC	NC		
Isophorone	μg	ND	ND	ND	NC	NC	NC		
2-Nitrophenol	μg	ND	ND	ND	NC	NC	NC		
2,4-Dimethylphenol	μg	ND	ND	ND	NC	NC	NC		
bis(2-Chloroethoxy) Methane	μg	ND	ND	ND	NC	NC	NC		
2,4-dichlorophenol	μg	ND	ND	ND	NC	NC	NC		
1,2,4-Trichlorobenzene	μg	0.75 J/J	ND	ND	NC	NC	NC		
Naphthalene	μg	18	6.2	0.97 J/J	NC	NC	NC		
4-Chloroaniline	μg	ND	ND	ND	NC	NC	NC		
Hexachlorobutadiene	μg	ND	ND	ND	NC	NC	NC		
4-Chloro-3-methylphenol	μg	ND	ND	ND	NC	NC	NC		
2-Methylnaphthalene	μв	5.4	ND	ND	NC	100.00%	NC		
Hexachlorocyclopentadiene	μв	ND	ND	ND	NC	NC	NC		
2,4,6-Trichlorophenol	μв	ND	ND	ND	NC	NC	NC		
2,4,5-Trichlorophenol	μg	ND	ND	ND	NC	NC	NC		
2-Chloronaphthalene	μg	ND	ND	ND	NC	NC	NC		
2-Nitroaniline	μg	ND	ND	ND	NC	NC	NC		
Dimethylphthalate	μg	ND	ND	NĐ	NC	NC	NC		
Acenaphthylene	μg	ND	ND	ND	NC	NC	NC		
2,6-Dinitrotoluene	μд	ND	ND	ND	NC	NC	NC NC		
3-Nitroaniline	μд	ND	ND	ND	NC	NC	NC		
Acenaphthene	μв	ND	ND	ND	NC	NC	NC		
2,4-Dinitrophenol	μg	ND	ND	ND	NC	NC	NC		
4-Nitrophenol	μg	ND	ND	ND	NC	NC	NC		
2,4-Dinitrotoluene	μg	ND	ND	ND	NC	NC	NC		
Dibenzofuran	μg	ND	ND	ND	NC	NC	NC		
Diethylphthalate	μg	ND	ND	. ND	NC	NC	NC		
Fluorene	μg	ND	ND	ND	NC	NC	NC NC		
-Chlorophenyl-phenyl Ether	μg	ND	ND	ND	NC	NC	NC		
4-Nitroaniline	μg	ND ND	ND	ND	NC NC	NC	NC		
1,6-Dinitro-2-methylphenol	μg	ND	ND	ND	NC NC	NC	NC		
N-Nitrosodiphenylamine	μg	ND ND	ND	ND	NC NC	NC NC	NC NC		
1-Bromophenyl-phenyl Ether	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC		
Hexachlorobenzene	μg	ND ND	ND	ND ND	NC NC	NC NC	NC NC		
Pentachlorophenol Phenanthrene	μg	ND ND	ND ND	ND ND	NC NC		NC NC		
Anthracene	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC		
di-n-Butylphthalate	μg	ND ND	ND	ND ND	NC NC	NC NC	NC NC		
luoranthene	μg μg	ND	ND	ND ND	NC NC	NC	NC NC		
Pyrene		ND ND	ND ND	ND	NC	NC	NC NC		
Sutylbenzylphthalate	μ <u>g</u>	ND	ND	ND	NC	NC	NC NC		
3,3'-Dichlorobenzidine	μg μg	ND	ND	ND	NC	NC	NC		
Chrysene		ND	ND	ND	NC	NC	NC		
Benzo(a)anthracene	μg μg	ND	ND	ND	NC	NC NC	NC		
pis(2-Ethylhexyl)phthalate	μg	2 J/J	ND	1.2 J/J	NC	NC	NC		
Di-n-Octylphthalate	μд	ND	ND	ND ND	NC	NC	NC		
Benzo(b)fluoranthene	μg	ND	ND	ND	NC	NC	NC		
Benzo(k)fluoranthene	μд	ND	ND	ND	NC	NC	NC		
Benzo(a)pyrene	μg	ND	ND	ND	NC	NC	NC		
indeno(1,2,3-c,d)pyrene	μg	ND	ND	ND	NC	NC NC	NC		
Dibenz(a,h)anthracene	μg	ND	ND	ND	NC	NC NC	NC		
Benzo(g,h,i)perylene	μg	ND	ND	ND	NC	NC	NC		
- ABITITY There	1	:	,::=,						
Fotol	1 115	EC 40	10.10	7 60	74.269	95.39%	84.82%		
Total	μg	56.40	10.10	2.60	74.26%	73.3770	07.0470		

Notes: _/ - Laboratory data qualifier /_ - Data validation qualifier

μg - Microgram

NC - Not calculated

ND · Non-detect

Destruction efficiency is not calculated where influent and/or effluent values are estimated.

Oualifiers: J · Result is estimated

1 The low destruction efficiency was not calculated because no SVOC compounds were detected in IN2.

American Chemical Service, Griffith, Indiana

		Round 2 - Sampled 5/22/02					
			Analytical Dat			truction Ef	
Compounds	Units	Influent IN1	Influent IN2	Effluent EF1	Low (%)	High (%)	Average (%)
Method TO-13							
Phenol	μg	ND	ND	ND	NC	NC	NC
bis(2-Chloroethyl)ether	μg	0.87 J/J	0.70 J/J	ND	NC	NC	NC NC
2-Chlorophenol	μg	ND	ND	ND	NC NC	NC	NC NC
1,3-Dichlorobenzene 1,4-Dichlorobenzene	μg	0.50 J/J	ND 4.4	ND ND	NC 100.00%	NC 100,000	NC 100.00%
1,2-Dichlorobenzene	μg	6.3	9.2	ND ND	100.00%	100.00%	100.00%
2-Methylphenol (o-Cresol)	μg	ND	ND	ND	NC	NC	NC
N-Nitroso-di-n-propylamine	μg	ND	ND	ND	NC NC	NC NC	NC NC
4-Methylphenol	μе	ND	ND	ND	NC	NC	NC NC
Hexachloroethane	μв	ND	ND	ND	NC	NC NC	NC
Nitrobenzene	μg	ND	ND	ND	NC	NC	NC
Isophorone	μg	ND	ND	ND	NC	NC	NC
2-Nitrophenol	μв	ND	ND	ND	NC	NC	NC
2,4-Dimethylphenol	μg	ND	ND	ND	NC	NC	NC
bis(2-Chloroethoxy) Methane	μg	ND	ND	ND	NC	NC	NC
2,4-dichlorophenol	μд	ND	ND	ND	NC	NC	NC
1,2,4-Trichlorobenzene	μg	1.6	1.2	ND	100.00%	100.00%	100.00%
Naphthalene	μg	29	23	ND	100.00%	100.00%	100.00%
4-Chloroaniline	μg	ND	ND	ND	NC_	NC	NC
Hexachlorobutadiene	μg	ND ND	ND ND	ND	NC	NC NC	NC NC
4-Chloro-3-methylphenol	μg	ND OA	ND 7.2	ND	NC 100,000	NC 100 000	NC 100,000
2-Methylnaphthalene	μg	9.4 ND	7.2 ND	ND ND	100.00% NC	100.00% NC	100.00% NC
Hexachlorocyclopentadiene 2,4,6-Trichlorophenol	μg	ND	ND	ND	NC	NC NC	NC NC
2,4,5-Trichlorophenol	μg	ND	ND	ND	NC	NC NC	NC
2-Chloronaphthalene	μg	ND	ND	ND	NC	NC NC	NC
2-Nitroaniline	μв	ND	ND	ND	NC	NC	NC
Dimethylphthalate	μg	ND	ND	ND	NC	NC	NC
Acenaphthylene	μg	ND	ND	ND	NC	NC	NC
2,6-Dinitrotoluene	μg	ND	ND	ND	NC	NC	NC
3-Nitroaniline	μg	ND	ND	ND	NC	NC	NC
Acenaphthene	μв	ND	ND	ND	NC	NC	NC
2,4-Dinitrophenol	μg	ND	ND	ND	NC	NC	NC
4-Nitrophenol	μg	ND	ND	ND	NC	NC	NC
2,4-Dinitrotoluene	1_ <u>4B</u> _	ND	ND	ND	NC	NC	NC
Dibenzofuran	μg	ND ND	ND	ND ND	NC NC	NC NC	NC NC
Diethylphthalate	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC
Fluorene 4-Chlorophenyl-phenyl Ether	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC
4-Nitroaniline	μg	ND	ND	ND ND	NC NC	NC NC	NC NC
4,6-Dinitro-2-methylphenol	μg μg	ND	ND	ND	NC	NC	NC NC
N-Nitrosodiphenylamine	μg	ND	ND	ND	NC NC	NC	NC
4-Bromophenyl-phenyl Ether	μg	ND	ND	ND	NC	NC	NC
Hexachlorobenzene	μд	ND	ND	ND	NC	NC	NC
Pentachlorophenol	μg	ND	ND	ND	NC	NC	NC
Phenanthrene	μв	ND	ND	ND	NC	NC	NC
Anthracene	μв	ND	ND	ND	NC	NC	NC
di-n-Butylphthalate	μg	0.85 J/J	0.81 J/J	ND	NC_	NC	NC
Fluoranthene	μg	ND	ND	ND	NC NC	NC NC	NC NC
Pyrene	μg	ND ND	ND ND	ND	NC NC	NC NC	NC NC
Butylbenzylphthalate 3,3'-Dichlorobenzidine	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC
Chrysene	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC
Benzo(a)anthracene	μg μg	ND	ND	ND	NC	NC	NC NC
bis(2-Ethylhexyl)phthalate	μg	ND	ND	ND	NC	NC	NC NC
Di-n-Octylphthalate	μg	ND	ND	ND	NC	NC	NC NC
Benzo(b)fluoranthene	μg	ND	ND	ND	NC	NC	NC
Benzo(k)fluoranthene	μg	ND	ND	ND	NC	NC	NC
Benzo(a)pyrene	μg	ND	ND	ND	NC	NC	NC
Indeno(1,2,3-c,d)pyrene	μg	ND	ND	ND	NC	NC	NC
Dibenz(a,h)anthracene	μg	ND	ND	ND	NC	NC	NC
Benzo(g,h,i)perylene	μg	ND	ND	ND	NC	NC	NC
. ,	↓ ↓ ↓				-		
Total	μе	59.30	45.00	ND	100.00%	100.00%	100.00%

Notes:
_/ - Laboratory data qualifier
/_ - Data validation qualifier
µg - Microgram

NC - Not calculated

ND - Non-detect

Destruction efficiency is not calculated where influent and/or effluent values are estimated

Ovalifiers: J - Result is estimated

1. The low destruction efficiency was not calculated because no SVOC

compounds were detected in IN2

American Chemical Service, Griffith, Indiana

		Round 3 - Sample						
			Analytical Dat			struction Ef		
Compounds	Units	Influent IN1	Influent IN2	Effluent EF1	Low (%)	High (%)	Average (%	
Method TO-13 Phenol		ND	ND	l MD	NC	I NC	l NG	
bis(2-Chloroethyl)ether	μg	ND ND	ND 1.2	ND ND	NC NC	NC NC	NC NC	
2-Chlorophenol	μg μg	ND	ND	ND ND	NC NC	NC NC	NC NC	
1,3-Dichlorobenzene	μg	ND	3.0	ND	NC NC	100.00%	NC NC	
1.4-Dichlorobenzene	μg	3.4	32	0.75 J/J	NC	NC	NC	
1,2-Dichlorobenzene	μg	6.1	56	1.2	80.33%	97.86%	89.09%	
2-Methylphenol (o-Cresol)	μg	ND	ND	ND	NC	NC	NC	
N-Nitroso-di-n-propylamine	μg	ND	ND	ND	NC	NC	NC	
4-Methylphenol	μg	ND	ND	ND	NC	NC	NC	
Hexachloroethane	μв	ND	ND	ND	NC	NC	NC	
Nitrobenzene	μg	ND	ND	ND	NC	NC	NC	
Isophorone	μg	ND	ND	ND	NC	NC	NC	
2-Nitrophenol	μg	ND	ND	ND	NC	NC	NC	
2,4-Dimethylphenol	μg	ND	ND_	ND	NC	NC	NC	
bis(2-Chloroethoxy) Methane	μg	ND	ND	ND	NC	NC	NC	
2,4-dichlorophenol	μе	ND .	ND	ND ND	NC	NC 100 000	NC NC	
1,2,4-Trichlorobenzene Naphthalene	μg	0.66 J/J 2.8	7.1	ND ND	NC 100.00%	100.00% 100.00%	NC 100.00%	
Naphthalene 4-Chloroaniline	μg	ND	ND	ND ND	NC	NC	NC	
Hexachlorobutadiene	μg μg	ND	ND	ND ND	NC	NC NC	NC NC	
4-Chloro-3-methylphenol	μg	ND	ND	ND	NC	NC NC	NC NC	
2-Methylnaphthalene	μв	1.4	15	ND	100.00%	100.00%	100.00%	
Hexachlorocyclopentadiene	μд	ND	ND	ND	NC	NC	NC	
2,4,6-Trichlorophenol	μв	ND	ND	ND	NC	NC	NC	
2,4,5-Trichlorophenol	μg	ND	ND	ND	NC	NC	NC	
2-Chloronaphthalene	μg	ND	ND	ND	NC	NC	NC	
2-Nitroaniline	μg	ND	ND	ND	NC	NC	NC	
Dimethylphthalate	μg	ND	ND	ND	NC	NC	NC	
Acenaphthylene	μg	ND	ND	ND	NC	NC	NC _	
2,6-Dinitrotoluene	μg	ND	ND	ND	NC	NC	NC	
3-Nitroaniline	μg	ND	ND	ND	NC	NC NC	NC NC	
Acenaphthene	μg	ND ND	ND	ND ND	NC NC	NC NC	NC	
2,4-Dinitrophenol	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC	
4-Nitrophenol 2,4-Dinitrotoluene	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC	
Dibenzofuran	μg μg	ND	ND	ND	NC	NC	NC NC	
Diethylphthalate	μв	0.37 J/J	0.43 J/J	0.34 J/J	NC NC	NC	NC	
Fluorene	μд	ND	ND	ND	NC	NC	NC	
4-Chlorophenyl-phenyl Ether	μg	ND	ND	ND	NC	NC	NC	
4-Nitroaniline	μв	ND	ND	ND	NC	NC	NC	
4.6-Dinitro-2-methylphenol	μд	ND	ND	ND	NC	NC	NC	
N-Nitrosodiphenylamine	μg	ND	ND	ND	NC	NC	NC	
1-Bromophenyl-phenyl Ether	μg	ND	ND	ND	NC	NC	NC	
Hexachlorobenzene	μg	ND	ND	ND	NC	NC	NC	
Pentachlorophenol	μg	ND	ND	ND	NC	NC	NC	
Phenanthrene	μg	ND	ND	ND ND	NC NC	NC NC	NC NC	
Anthracene	μg	ND ND	ND 14 W	ND ND	NC NC	NC NC	NC NC	
li-n-Butylphthalate Tuoranthene	μg μg	ND ND	1.4 J/J ND	ND ND	NC NC	NC NC	NC NC	
Pyrene		ND ND	ND ND	ND ND	NC	NC	NC NC	
Butylbenzylphthalate	μg	ND ND	ND	ND	NC	NC NC	NC	
3,3'-Dichlorobenzidine	μg	ND	ND	ND ND	NC	NC	NC NC	
Chrysene	μg	ND	ND	ND	NC	NC	NC	
Benzo(a)anthracene	μg	ND	ND	ND	NC	NC	NC _	
ois(2-Ethylhexyl)phthalate	μg	2.0 J/J	1.9 J/J	6.1	NC	NC	NC	
Di-n-Octylphthalate	μg	ND	ND	ND	NC	NC	NC	
Benzo(b)fluoranthene	μg	ND	ND	ND	NC	NC	NC	
Benzo(k)fluoranthene	μg	ND	ND	ND	NC	NC	NC	
Benzo(a)pyrene	μg	ND	ND	ND	NC	NC	NC	
ndeno(1,2,3-c,d)pyrene	μg	ND	ND	ND	NC	NC	NC	
Dibenz(a,h)anthracene	μg	ND ND	ND	ND	NC	NC NC	NC NC	
Benzo(g,h,i)perylene	μg	ND	ND	ND	NC	NC	NC	
	+	· · · · · · · · · · · · · · · · · · ·						
l'otal	μg	13.70	141.30	7.30	46.72%	94.83%	70.77%	

Oualiflers: J - Result is estimated

Notes:
_/ - Laboratory data qualifier
/_ - Data validation qualifier

JB - Analyte is detected in the method blank resulting in potential bias high Reported concentration is estimated

μg - Microgram

NC - Not calculated

ND - Non-detect

Destruction efficiency is not calculated where influent and/or effluent values are estimated.

1. The low destruction efficiency was not calculated because no SVOC

compounds were detected in IN2

American Chemical Service, Griffith, Indiana

Γ	_		Re	ound 4 - Samples	1 6/28/02		
			Analytical Data			struction Eff	iciency
Compounds	Units	Influent IN1	Influent IN2	Effluent EF1	Low (%)	High (%)	Average (%)
Method TO-13				pi		· · ·	·
Phenol	μg	ND	ND	ND	NC_	NC	NC NC
bis(2-Chloroethyl)ether	μg	1.7	ND	ND	NC	100.00%	NC
2-Chlorophenol	μg	ND	ND	ND	NC NC	NC 100.007	NC NC
1,3-Dichlorobenzene	μg	2.7	ND	ND	NC NC	100.00%	NC NC
1,4-Dichlorobenzene 1,2-Dichlorobenzene	μg	27 46	ND ND	1.5 2.3	NC NC	94.44% 95.00%	NC NC
2-Methylphenol (o-Cresol)	μg	ND ND	ND	ND ND	NC	NC	NC NC
N-Nitroso-di-n-propylamine	Pg Bu	ND	ND	ND	NC	NC NC	NC
4-Methylphenol	μg	ND	ND	ND	NC	NC	NC
Hexachloroethane	μg	ND	ND	ND	NC	NC	NC
Nitrobenzene	μg	ND	ND	ND	NC	NC	NC
Isophorone	μg	ND	ND	ND	NC	NC	NC
2-Nitrophenol	μg	ND	ND	ND	NC	NC	NC
2,4-Dimethylphenol	μg	ND	ND_	ND	NC	NC	NC
bis(2-Chloroethoxy) Methane	μg	ND	ND	ND	NC	NC	NC
2,4-dichlorophenol	μg	ND	ND	ND	NC	NC	NC
1,2,4-Trichlorobenzene	μg	6.2	ND	0.38 J/J	NC	NC NC	NC NC
Naphthalene	μg	13	ND ND	0.41 J/J	NCNC	NC NC	NC_
4-Chloroaniline Hexachlorobutadiene	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC
Hexachlorobutadiene 4-Chloro-3-methylphenol	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC
2-Methylnaphthalene	μg	6.9	ND ND	ND ND	NC NC	100.00%	NC NC
Hexachlorocyclopentadiene	μg μg	ND	ND	ND	NC NC	NC	NC
2,4,6-Trichlorophenol	μg	ND	ND	ND	NC	NC	NC
2.4.5-Trichlorophenol	μg	ND	ND	ND	NC	NC	NC
2-Chloronaphthalene	μв	ND	ND	ND	NC	NC	NC
2-Nitroaniline	μв	ND	ND	ND	NC	NC	NC
Dimethylphthalate	μg	ND	ND	ND	NC	NC	NC
Acenaphthylene	μg	ND	ND	ND	NC	NC	NC NC
2,6-Dinitrotoluene	μg	ND	ND	ND	NC	NC	NC
3-Nitroaniline	μg	ND	ND	ND	NC	NC	NC
Acenaphthene	μв	ND	ND	ND	NC NC	NC NC	NC
2,4-Dinitrophenol	μg	ND	ND	ND ND	NC NC	NC NC	NC NC
4-Nitrophenol 2.4-Dinitrotoluene	μg	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC
Dibenzofuran	μg	ND	ND ND	ND	NC NC	NC NC	NC NC
Diethylphthalate	μg μg	0.52 J/JB	ND	0.47 J/JB	NC	NC NC	NC NC
Fluorene	μв	ND	ND	ND	NC	NC NC	NC
4-Chlorophenyl-phenyl Ether	μg	ND	ND	ND	NC	NC	NC
4-Nitroaniline	μg	ND	ND	ND	NC	NC	NC
4,6-Dinitro-2-methylphenol	μд	ND	ND	ND	NC	NC	NC
N-Nitrosodiphenylamine	μg	ND	ND	ND	NC	NC	NC
4-Bromophenyl-phenyl Ether	μg	ND	ND	ND	NC	NC	NC
Hexachlorobenzene	μg	ND	ND	ND	NC	NC	NC
Pentachlorophenol	μg	ND ND	ND	ND	NC NC	NC NC	NC
Phenanthrene	⊢.μg 	ND ND	ND ND	ND ND	NC NC	NC NC	NC NC
Anthracene di-n-Butylphthalate	μg	1.2 J/J	ND ND	ND ND	NC NC	NC NC	NC NC
Fluoranthene	μg μg	ND ND	ND ND	ND	NC NC	NC NC	NC NC
Pyrene	μg	ND	ND	ND ND	NC	NC NC	NC NC
Butylbenzylphthalate	μg	ND	ND	ND	NC	NC	NC
3,3'-Dichlorobenzidine	μв	ND	ND	ND	NC	NC	NC
Chrysene	μg	ND	ND	ND	NC	NC	NC _
Benzo(a)anthracene	μg	ND	ND	ND	NC	NC	NC
bis(2-Ethylhexyl)phthalate	μg	ND	ND	ND	NC	NC	NC
Di-n-Octylphthalate	μg	ND	ND	ND	NC	NC	NC
Benzo(b)fluoranthene	μg	ND	ND	ND	NC	NC NC	NC
Benzo(k)fluoranthene	μg	ND	ND	ND	NC	NC	NC
Benzo(a)pyrene	μg	ND	ND	ND	NC	NC	NC
Indeno(1,2,3-c,d)pyrene	μg	ND	ND	ND	NC NC	NC NC	NC NC
Dibenz(a,h)anthracene	μg	ND ND	ND	ND	NC NC	NC NC	NC NC
Benzo(g,h,i)perylene	μg	ND	ND	ND	NC	NC	NC
Total	μg	103.50	ND	3.80	NC	96.33%	NC

Notes: _/ - Laboratory data qualifier /_ - Data validation qualifier

Oualifiers: J · Result is estimated

μg - Microgram

IB - Analyte is detected in the method blank resulting in potential

bias high. Reported concentration is estimated

NC - Not calculated

ND - Non-detect

Destruction efficiency is not calculated where influent and/or effluent values are estimated

1 The low destruction efficiency was not calculated because no SVOC

compounds were detected in IN2

Table 4.1
Water Levels Inside Barrier Wall - Second Quarter 2002
American Chemical Service NPL Site
Griffith, Indiana

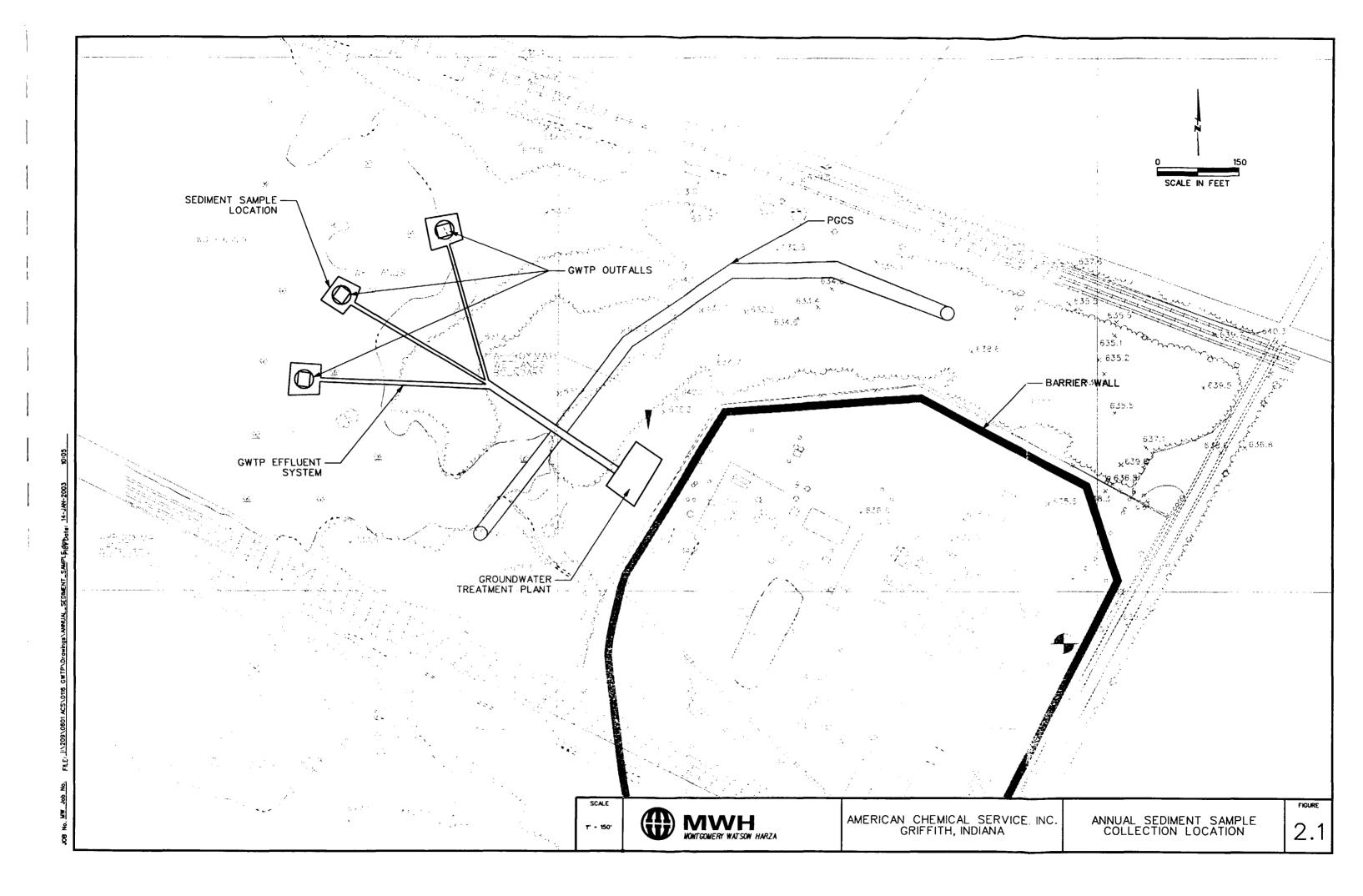
Data	On-Site Area									
Date	P-29	P-31	P-32	P-36	P-49	P-106	P-108			
5-Apr-02	633.7	635.4	634.4	635.0	632.3	633.0	634.3			
12-Apr-02	633.7	635.5	634.5	635.0	632.6	633.1	634.4			
19-Apr-02	633.8	635.5	634.4	635.0	632.6	632.8	634.1			
26-Apr-02	633.7	636.0	634.4	633.2	634.1	634.9	636.3			
3-May-02	633.6	636.4	634.3	631.3	635.5	636.9	638.4			
10-May-02	633.9	636.7	634.5	631.5	636.1	637.2	638.8			
17-May-02	634.7	637.7	635.2	632.6	637.3	637.9	639.4			
24-May-02	634.4	637.0	634.8	632.2	636.5	637.6	639.2			
31-May-02	634.4	636.8	634.5	632.2	636.6	637.9	638.5			
7-Jun-02	634.1	636.5	634.3	631.8	636.1	637.5	638.2			
14-Jun-02	633.3	635.9	634.0	631.3	635.5	636.9	637.9			
21-Jun-02	632.6	635.3	633.7	630.9	634.8	636.4	637.7			
28-Jun-02	633.2	635.0	633.8	630.7	634.8	636.1	638.0			

D-4-	Off-Site Area								
Date	P-96	P-110	P-112	P-113	P-114	P-116	P-118		
5-Apr-02	620.9	628.8	625.7	628.5	629.5	628.9	627.6		
12-Apr-02	622.6	629.0	626.9	628.4	629.3	628.8	627.4		
19-Apr-02	620.9	628.5	625.7	628.2	629.1	628.6	627.5		
26-Apr-02	620.9	628.3	626.1	628.2	629.0	628.5	627.3		
3-May-02	620.9	628.2	626.4	628.0	628.9	628.4	627.0		
10-May-02	620.9	628.3	627.0	628.0	628.9	628.5	627.0		
17-May-02	620.9	628.7	627.3	627.9	628.9	628.4	626.3		
24-May-02	620.9	628.4	624.4	627.8	628.9	627.9	627.0		
31-May-02	620.9	628.5	624.8	628.7	628.9	628.3	626.9		
7-Jun-02	620.9	628.1	624.2	627.1	627.5	627.9	626.7		
14-Jun-02	620.9	627.9	624.1	627.3	628.0	627.8	626.6		
21-Jun-02	620.9	627.8	624.1	627.4	628.5	627.7	626.5		
28-Jun-02	620.9	627.7	624.0	627.3	628.4	627.8	626.6		

Notes:

All water level elevations are in feet AMSL





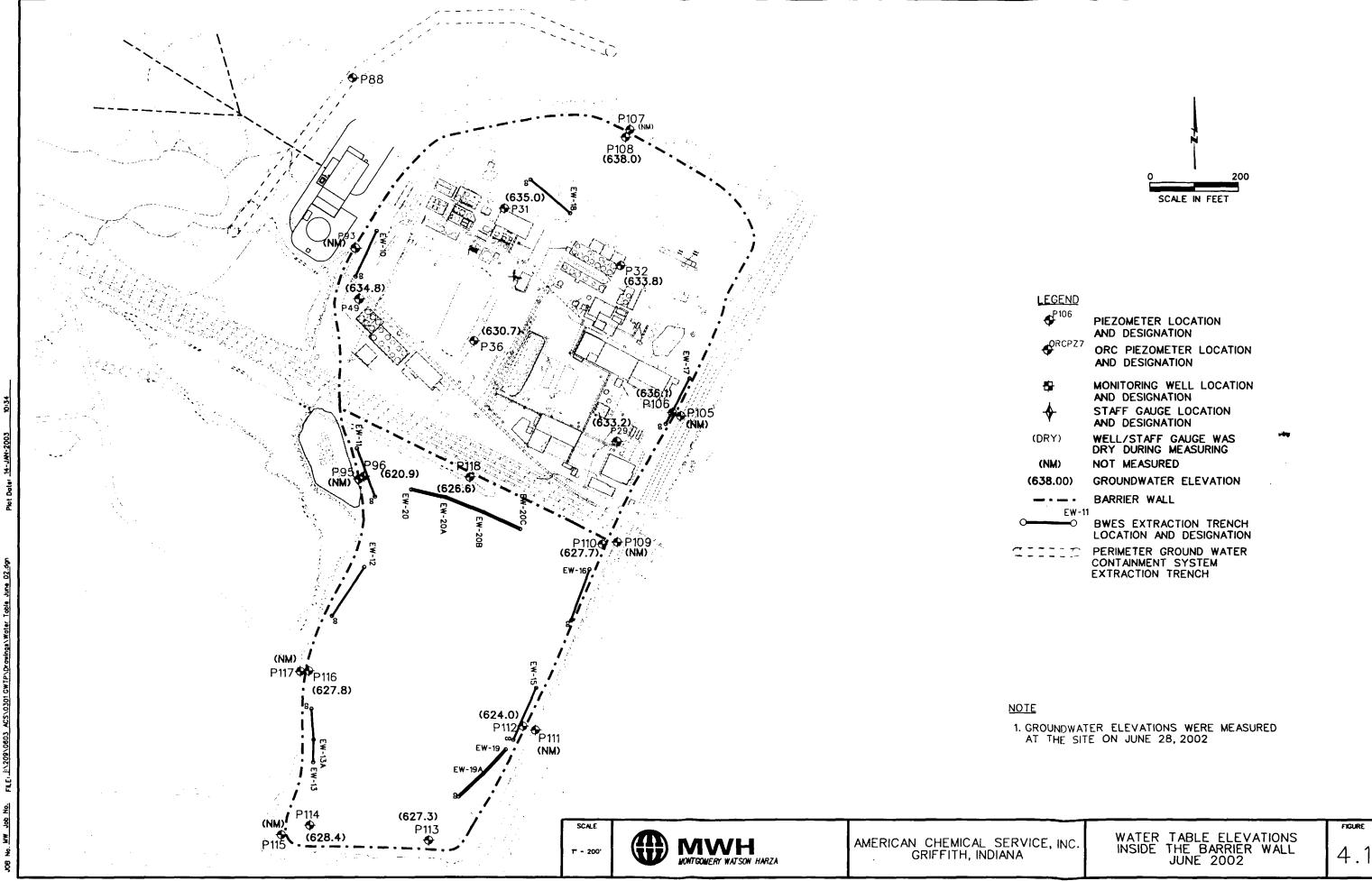
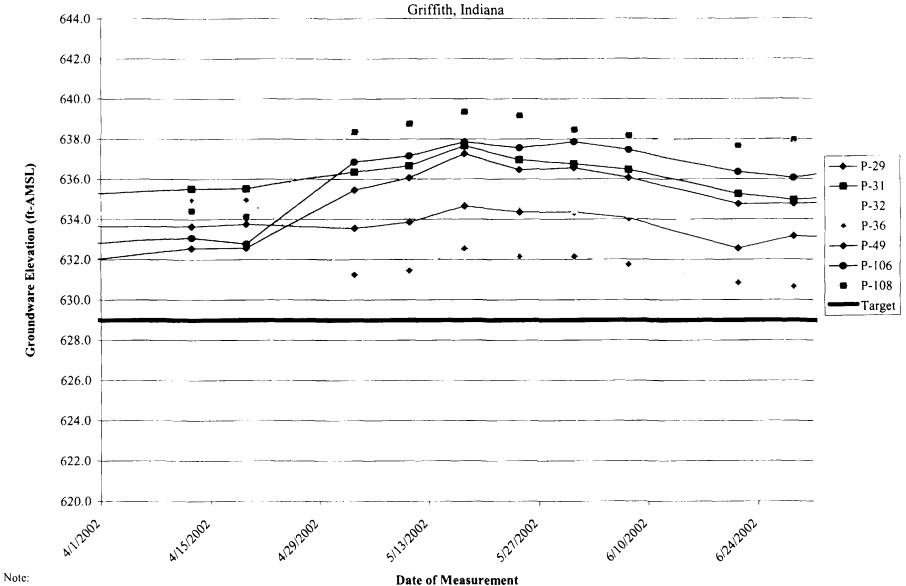


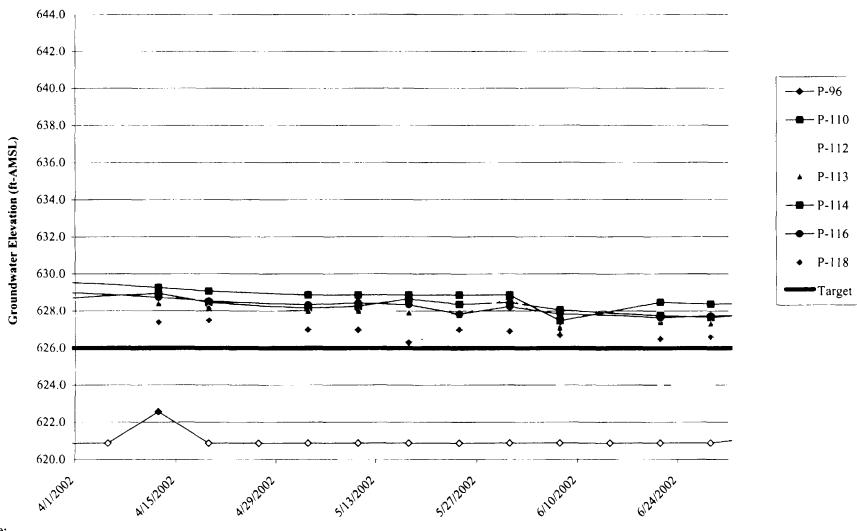
Figure 4.2
Water Level Trends Inside Barrier Wall (On-Site Area)
ACS NPL Site



Hollow points represent dry piezometers (data used for graphing purposes only).

TMK/RHS
J:/209/0603/0301/6030301a019.xls/On-Site Chart for Report

Figure 4.3
Water Level Trends Inside Barrier Wall (Off-Site Area)
ACS NPL Site
Griffith, Indiana



Note:

Hollow points represent dry piezometers (data used for graphing purposes only).

Date of Measurement

TMK/RHS

J:/209/0603/0301/6030301a019.xls/Off-Site Chart for Report

APPENDIX A EFFLUENT ANALYTICAL DATA

April 22, 2002 Compliance Sample Laboratory Results

SW-846

1-CC

CLASSICAL CHEMISTRY ANALYSES DATA SHEET

EPA SAMPLE NO.

Lar Name: CompuC		CompuChem		Contract:			EFFLUENT			
	Code:	LIBRTY	Case No.: NRAS N							
SDG	No.:	RC1024								
Mat	ix (so	oil/water):	WATER		La	ab Sam	ple I	D: RC1024-1		
Dat	e Recei	ved: 4/23/0	2		*	Solid	ls: 0	.00		
		Concen	tration Un	its (mg/L or mg/kg o	lry weigh	t):	mg	/L		
		PARAMETI	ER _	CONCENTRATI	ON C	Q	М	DATE ANALYZED]	
		TSS	·	1	.00 ט			4/23/02	Ī	
		pН		7	. 96			4/23/02]	

Comments:

CHEMICAL & ENVIRONMENTAL TECHNOLOGY, INC.

SUUSB

ENVIRONMENTAL ANALYTICAL SERVICES

FINAL REPORT OF ANALYSES

COMPUCHEM
Actn: DIANE BYRD
501 MADISON AVENUE
CARY, NC 27513-

REPORT DATE: 05/14/02

ACS-89

SAMPLE NUMBER- 196307 SAMPLE ID- EFFLUENT

DATE SAMPLED- 04/22/02

DATE RECEIVED- 04/23/02 SAMPLER- NOT SPECIFIED

TIME RECEIVED- 1500

DELIVERED BY- CHRIS BRAND

DEDITION DI CHILD BIVE

Page 1 of 1

PROJECT NAME : ACS-89

ANALYSIS

ANALYSIS

METHOD DATE

BY RESULT UNITS

PQL

SAMPLE MATRIX- WW TIME SAMPLED- 1400

RECEIVED BY- RCB

BIOCHEMICAL OXYGEN DEMAND

EPA 405.1 04/24/02 LEB

<2 mg/L

2

PQL - Practical Quantitation Limit

Results followed by the letter 3 are estimated concentrations.

NC DENR CERTIFICATIONS: DWQ - 96; PUBLIC WATER SUPPLY - 37724

LABORATORY DIRECTOR

15241

SW-846 METALS

ı

INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

EFFLUENT

	COMPUCHEM		Contr	act:			_	
ode:	LIBRTY	Case No.:		SAS No.:		<u>.</u>	SDG 1	No.: RC1024
x (soi	l/water):	WATER		Lab Sample II) :	RC102	4-1	
l (low/	med): LC	n u		Date Receive	4 :	04/23	/02	
(10.17	<u></u>			2200 110021701	•••	04/25/	02_	
lids:	0.0							
		Gamaantwati an	. 17mita (22m/7)					-
		Concentration	onits (ug/L o	or mg/kg dry weig	nt)	•	UG/	<u> 4</u>
		CAS No.	Analyte	Concentration	c	Q	М]
		17400 00 5		1			 _	iub
		7429-90-5	Aluminum	88.5		<u> </u>		
		7440-36-0	Antimony	4.1		<u> </u>	P	<u> โ</u> นช
		7440-38-2	Arsenic	2.0	ט [<u> </u>	P	-
		7440-39-3	Barium	70.1	1	<u> </u> 	P	-
		7440-41-7	Beryllium	0.20		<u> </u>	P	-
		7440-70-2	Calcium	68700	10	<u> </u>	P	-1
		7440-47-3	Chromium	0.50	 77		P P	1
		7440-47-3	Cobalt	1.2		<u> </u>	l P	-
		7440-50-8	Copper	1.0			P	1
		7439-89-6	Iron	9.5		<u> </u>	P	4
		7439-92-1	Lead	2.3		<u> </u>	P	iub
		7439-95-4	Magnesium	29500		<u> </u>	I P	1000
		7439-96-5	Manganese	14.4		<u> </u>	P	1
		7439-97-6	Mercury	0.10	<u> </u>		CV	i
		7440-02-0	Nickel	5.1			P	i
		7440-09-7	Potassium	12500	<u> </u>		P	j
		7782-49-2	Selenium	2.1	U		P	Ī
		7440-22-4	Silver	0.50			P	Ĭ
		7440-23-5	Sodium	87700	1		P	Ī
		7440-28-0	Thallium	2.2	U	}	P	Ī
		7440-62-2	Vanadium	3.8		i	P	Ī
			Zinc	7.1	1 -		$\overline{}$	inb

Color Before:	COLORLESS	Clarity Before:	CLEAR	Texture:	
olor After:	COLORLESS	Clarity After:	CLEAR	Artifacts:	
Comments:					
					8

Q

FORM 1 VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM Method: 3260B EFFLUENT

ab Code: LIBRTY Case No.: SAS No.: SDG No.: RC1024

Matrix: (soil/water) WATER Lab Sample ID: RC1024-1

ample wt/vol: 25 (g/ml) ML Lab File ID: RC1024-1B73

revel: (low/med) LOW Date Received: 04/23/02

Moisture: not dec. Date Analyzed: 04/25/02

C Column: ZB-624 ID: 0.32 (mm) Dilution Factor: 1.0

Soil Extract Volume: (uL) Soil Aliquot Volume: (uI

CAS NO. COMPOUND CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.	COMPOUND (ag/ II of ag	,, kg, 00, 11		Q	
74-87-3	Chloromethane		0.1	J	
	Vinyl Chloride	1 (0.5	บ	
	Bromomethane		0.5		
	Chloroethane	-).5	ט ו	
75-35-4	1,1-Dichloroethene	-) c	0.5	บ	
	Carbon disulfide	·)	0.5	ULBL	2110
67-64-1		- [1	JB 3	BUB
	Methylene Chloride	-	0.1	J	
	trans-1,2-Dichloroethene		0.5		
	1,1-Dichloroethane	·	0.5	ប 🏻	
	cis-1,2-Dichloroethene	- (0.5	U	
	2-butanone	- }	3		
67-66-3	Chloroform	-	0.5	U	
71-55-6	1,1,1-Trichloroethane	-	0.5	บ	
	Carbon Tetrachloride	-) (0.5	ט	
71-43-2	Benzene		0.5	U	
107-06-2	1,2-Dichloroethane	-1 (0.51	υ	
79-01-6	Trichloroethene	- (0.5	ַ ט (
78-87-5	1,2-Dichloropropane	-)	0.5	ט ו	
	Bromodichloromethane	-) (0.5	U	
10061-01-5	cis-1,3-Dichloropropene	-	0.5	ט	
108-10-1	4-Methyl-2-pentanone	-	3	U	
108-88-3	Toluene	0	.06	J	
10061-02-6	trans-1,3-Dichloropropene	-	0.5	U	
79-00-5	1,1,2-Trichloroethane	-	0.5	υ 🚽	
127-18-4	Tetrachloroethene	-	.05	JB 0.5	UB
	2-hexanone	-	3	U	
124-48-1	Dibromochloromethane	-	0.5	ប	
108-90-7	Chlorobenzene	-	0.5	U	}
100-41-4	Ethylbenzene	-	0.5	U	1
108-38-3	m,p-Xylene	_	1	U	(
95-47-6	o-Xylene	-	0.5	ប	1
	Styrene		0.5	ប	
l	EODM T VOX	_	!	l	1

FORM I VOA

SP 11

FORM 1 VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EFFLUENT

Method: 8260B Lab Name: COMPUCHEM SDG No.: RC1024 ab Code: LIBRTY Case No.: SAS No.: Matrix: (soil/water) WATER Lab Sample ID: RC1024-1 Lab File ID: RC1024-1B73 ample wt/vol: 25 (q/ml) ML Level: (low/med) LOW Date Received: 04/23/02 Moisture: not dec. Date Analyzed: 04/25/02 GC Column: ZB-624 ID: 0.32 (mm) Dilution Factor: 1.0 Soil Aliquot Volume: ____(uL oil Extract Volume: ____ (uL) CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/L 0 75-25-2-----Bromoform 0.5 U 79-34-5----1,1,2,2-Tetrachloroethane 0.5 U 541-73-1----1,3-Dichlorobenzene____ 0.5 U 106-46-7----1,4-Dichlorobenzene 0.5|U 95-50-1-----1,2-Dichlorobenzene 0.5 U 120-82-1----1,2,4-Trichlorobenzene 0.5 UUJ 540-59-0-----1,2-Dichloroethene (total) 0.5 U 1330-20-7-----Xylene (total) 0.5 U

FORM I VOA

6 12

FORM 1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EFFLUENT

Method: 8270C Lab Name: COMPUCHEM

Lab Code: LIBRTY Case No.: SAS No.: SDG No.: RC1024

Matrix: (soil/water) WATER

Lab Sample ID: RC1024-1

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: RC1024-1RA64

Level: (low/med) LOW

Date Received: 04/23/02

% Moisture: decanted: (Y/N)

Date Extracted: 04/29/02

Concentrated Extract Volume: 1000(uL)

Date Analyzed: 04/29/02

Injection Volume: 1.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _

CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/L

0

111-44-4Bis(2-chloroethyl)ether 106-44-54-Methylphenol 78-59-1Isophorone	9.6 10 10	U U	
78-59-1Isophorone 117-81-7bis(2-ethylhexyl)Phthalate	0.91	JB 6	u

FORM I SV

8270C

FORM 1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EFFLUENT

Lab Name: COMPUCHEM	Method:	SIM		
Lab Code: LIBRTY	Case No.: SAS No.	: SDG	No.:	RC1024
Matrix: (soil/water)	WATER	Lab Sample ID:	RC10	24-1
Sample wt/vol:	1000 (g/mL) ML	Lab File ID:	RC10	24-1A64
Level: (low/med)	LOW	Date Received:	: 04/2	3/02
k Moisture:	decanted: (Y/N)	Date Extracted	1:04/2	9/01
Concentrated Extract	Volume: 1000(uL)	Date Analyzed:	05/0	2/02
Injection Volume:	1.0(uL)	Dilution Facto	or: 1.	0
GPC Cleanup: (Y/N)	N pH:			
CAS NO.		NTRATION UNITS: or ug/Kg) UG/I	-	Q
87-86-5	Pentachlorophenol		1	U

FORM I SV

13

EFFLUENT

Lab Name: COMPUCHEM	Contract: 8082	
ab Code: LIBRTY Case No.:	SAS No.: SDG	No.: RC1024
Matrix: (soil/water) WATER	Lab Sample ID:	RC1024-1

ample wt/vol: 1000 (g/mL) ML Lab File ID:

% Moisture: decanted: (Y/N) Date Received: 04/23/02

xtraction: (SepF/Cont/Sonc) SEPF Date Extracted:04/25/02

Concentrated Extract Volume: 5000(uL) Date Analyzed: 04/26/02

njection Volume: 2.0(uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: ___ Sulfur Cleanup: (Y/N) N

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q 12674-11-2----Aroclor-1016 0.50 U 1.0 U 11104-28-2----Aroclor-1221 11141-16-5-----Aroclor-1232 0.50 U 0.50 53469-21-9-----Aroclor-1242 12672-29-6-----Aroclor-1248 0.50 11097-69-1----Aroclor-1254 0.50 ĺυ 11096-82-5----Aroclor-1260 0.50 U

CONCENTRATION UNITS:

Sold

May 9, 2002 Compliance Sample Laboratory Results

SW-846

1-CC

CLASSICAL CHEMISTRY ANALYSES DATA SHEET

EPA SAMPLE NO.

							EFFLU	ENT
ab lame:	CompuChem		Contract:			¹		
ab Code:	LIBRTY	Case	No.:		_	NRA	S No.:	
;DG 40.:	RE1024	_						
fatrix (soi	1/water):	WATER		La	b Sam	ple I	D: RE1024-1	
)at Receiv	red: 5/17/02			*	Solid	s: 0	.00	
	Concenti	ration Units (mg/	L or mg/kg dry	weight	:):	mg/	/L	
							DATE	
	PARAMETER		CONCENTRATION	С	Q	м	ANALYZED	
	TSS		1.00	บ			5/21/02	
	рH		7.83				5/17/02	

rayor

Comments:						

FINAL REPORT OF ANALYSES

COMPUCHEM

Attn: DIANE BYRD 501 MADISON AVENUE CARY, NC 27513-

REPORT DATE: 05/29/02

SAMPLE NUMBER- 197124 SAMPLE ID- EFFLUENT

DATE SAMPLED- 05/16/02

DATE RECEIVED- 05/17/02 SAMPLER- NOT SPECIFIED

TIME RECEIVED- 1350

DELIVERED BY- JAMES FELDHAUS

Page 1 of 1

PROJECT NAME : ACS 89

ANALYSIS

ANALYSIS

METHOD

DATE

BY RESULT UNITS

PQL

SAMPLE MATRIX- WW

RECEIVED BY- ALT

TIME SAMPLED- 1430

BIOCHEMICAL OXYGEN DEMAND

EPA 405.1 05/17/02 LEB

<2 mg/L

2

PQL = Practical Quantitation Limit

Results followed by the letter J are estimated concentrations.

NC DENR CERTIFICATIONS: DWQ - 96; PUBLIC WATER SUPPLY - 37724

LABORATORY DIRECTOR

710 010 010 WY 00 01 70/00/00

SW846 METALS

INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

al Name: COMPUCHEM		Contra	ict:			EFFLUENT
				_ :		
Lab Code: LIBRTY	Case No.:	SAS	No.:	SDG	No.:	RE1024
Mai ix (soil/water):	WATER		Lab Sample I	D: <u>RE10</u>	24-1	<u>_</u>
evel (low/med): LO	W		Date Receive	d: <u>5/</u> 17	/02	
lids: 0.0						•
	Concentrati	on Units (ug/L	or mg/kg dry	weight):	UG/I	<u>L</u>
	CAS No.	Analyte	Concentration	C Q	М	
<u> </u>	429-90-5	Aluminum	120		P	iuß
7	440-38-2	Arsenic	4.2	ן ט	P	1
7	440-39-3	Barium	75.5		P	Ī
<u> </u>	440-36-0	Antimony	3.8	B	P	iub
	440-41-7	Beryllium	0.21	В	P	IUB
7	440-43-9	Cadmium	0.30	ן ט	P	1
7	440-70-2	Calcium	64000		P	
[7	440-47-3	Chromium		B	P	iub
<u>\</u>	440-48-4	Cobalt		B	P	
!	440-50-8	Copper		B	P	
<u></u>	439-89-6	Iron		U	P	
<u>'</u>	439-92-1	Lead		ן ט ן	P	
]	439-95-4	Magnesium	28200	18 (P	
'	439-96-5	Manganese		B U	P	
<u>'</u>	439-97-6	Mercury		10 1	CV	 -
<u></u>	440-02-0	Nickel Potassium	9.9	<u> </u>	P	<u> </u>
<u></u>	782-49-2	Selenium	<u>'</u>	 U	I P	<u>t</u> 1
<u>'</u>	440-22-4	Silver		[U]	P	
<u></u>	440-23-5	Sodium	165000	 	l P	<u>!</u>
<u> </u>	440-28-0	Thallium	<u>' </u>	ו טו	P	<u>'</u>
<u></u>	440-62-2	Vanadium	<u> </u>	B	P	UB
 	440-66-6	Zinc	<u> </u>	וט ו	P	! W.D.
<u>. </u>		<u>'</u>	'	<u>· · · · · · · · · · · · · · · · · · · </u>		<u>:</u>
						rolyon
Color Before: COLOR	LESS Cla	arity Before:	CLEAR	Textur	e:	
Color After: COLOR	LESS Cla	arity After:	CLEAR	Artifa	cts:	·
Comments:						

FORM 1 VOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: COMPUCHEM Method: 8260B EFFLUENT

I b Code: LIBRTY Case No.: SAS No.: SDG No.: RE1024

Matrix: (soil/water) WATER Lab Sample ID: RE1024-1

Sample wt/vol: 25 (g/ml) ML Lab File ID: RE1024-1B71

: evel: (low/med) LOW Date Received: 05/17/02

% Moisture: not dec. Date Analyzed: 05/29/02

C Column: SPB-624 ID: 0.32 (mm) Dilution Factor: 1.0

Soil Extract Volume: (uL) Soil Aliquot Volume: (uL

CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/L O 74-87-3-----Chloromethane 0.5 U 75-01-4-----Vinyl Chloride 0.5 U 0.5 U 0.5 U 0.5 U 74-83-9-----Bromomethane 75-00-3-----Chloroethane 75-35-4----1,1-Dichloroethene 75-15-0-----Carbon disulfide 0.5 UU 67-64-1-----Acetone 75-09-2-----Methylene Chloride 156-60-5----trans-1,2-Dichloroethene 0.5 U 0.5 U 0.5 U 75-34-3-----1,1-Dichloroethane 156-59-2----cis-1,2-Dichloroethene 78-93-3----2-butanone 3 U 0.5 U 67-66-3-----Chloroform 71-55-6-----1,1,1-Trichloroethane 0.5] ប 56-23-5-----Carbon Tetrachloride 0.5 U 0.5 U 71-43-2-----Benzene 107-06-2----1,2-Dichloroethane 0.5 U 79-01-6-----Trichloroethene 0.5 U 78-87-5-----1,2-Dichloropropane 0.5 U 0.5|0 75-27-4-----Bromodichloromethane U 2.0 10061-01-5----cis-1,3-Dichloropropene 108-10-1-----4-Methyl-2-pentanone 108-88-3-----Toluene 0.2 JB 0.5 WB 10061-02-6----trans-1,3-Dichloropropene 0.5 U 0.5 U 79-00-5-----1,1,2-Trichloroethane 0.5 UUJ 127-18-4-----Tetrachloroethene 591-78-6----2-hexanone 124-48-1-----Dibromochloromethane 0.5|U 108-90-7-----Chlorobenzene 0.04 J 100-41-4-----Ethylbenzene 0.5 U 108-38-3----m,p-Xylene 1 | U 95-47-6----o-Xylene 0.5 U 100-42-5-----Styrene 0.510

FORM I VOA

FORM 1 VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EFFLUENT Lab Name: COMPUCHEM Method: 8260B lab Code: LIBRTY Case No.: SAS No.: SDG No.: RE1024 Matrix: (soil/water) WATER Lab Sample ID: RE1024-1 Lample wt/vol: 25 (g/ml) ML RE1024-1B71 Lab File ID: evel: (low/med) LOW Date Received: 05/17/02 * Moisture: not dec. Date Analyzed: 05/29/02 Dilution Factor: 1.0 C Column: SPB-624 ID: 0.32 (mm) Soil Aliquot Volume: ____(uL Soil Extract Volume: ____(uL) CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q 75-25-2-----Bromoform 0.5 U 79-34-5-----1,1,2,2-Tetrachloroethane 0.5 U 541-73-1----1,3-Dichlorobenzene 0.5 0 106-46-7-----1,4-Dichlorobenzene 0.5 U 95-50-1----1,2-Dichlorobenzene 0.5 U 120-82-1----1,2,4-Trichlorobenzene 0.5 U 540-59-0-----1,2-Dichloroethene (total) 0.5 U 1330-20-7-----Xylene (total) 0.5 U

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FORM I VOA

FORM 1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EFFLUENT

Lab Name: COMPUCHEM	Method:	EFFLUENT
ab Code: LIBRTY	Case No.: SAS No	.: SDG No.: RE1024
Matrix: (soil/water)	WATER	Lab Sample ID: RE1024-1
Sample wt/vol:	1000 (g/mL) ML	Lab File ID: RE1024-1A66
evel: (low/med)	LOW	Date Received: 05/17/02
% Moisture:	decanted: (Y/N)	Date Extracted:05/20/02
Concentrated Extract	Volume: 1000(uL)	Date Analyzed: 05/21/02
Injection Volume:	1.0(uL)	Dilution Factor: 1.0
GPC Cleanup: (Y/N)	N pH:	
CAS NO.		ENTRATION UNITS: L or ug/Kg) UG/L Q
•		i l

,6402

8270C

FORM 1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EFFLUENT

Q

_ab Code: LIBRTY Case No.: SAS No.: SDG No.: RE1024 Lab Sample ID: RE1024-1 Matrix: (soil/water) WATER Sample wt/vol: 1000 (g/mL) ML

Lab File ID: RE1024-1A70

Date Received: 05/17/02 Level: (low/med) LOW

% Moisture: _____ decanted: (Y/N)___ Date Extracted:05/20/02

Concentrated Extract Volume: 1000(uL) Date Analyzed: 05/23/02

Dilution Factor: 1.0 Injection Volume: 1.0(uL)

GPC Cleanup: (Y/N) N pH:____

Lab Name: COMPUCHEM

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L COMPOUND CAS NO.

87-86-5-----Pentachlorophenol____ U 1

Method:

1D GC EXTRACTABLE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

EFFLUENT

Lab Name: COMPUCHEM	Contract: 8082
La Code: LIBRTY Case No.:	SAS No.: SDG No.: RE1024
Matrix: (soil/water) WATER	Lab Sample ID: RE1024-1
Sample wt/vol: 1000 (g/mL) ML	Lab File ID:
% Moisture: decanted: (Y/N)	Date Received: 05/17/02
Extraction: (SepF/Cont/Sonc) SEPF	Date Extracted:05/17/02
C ncentrated Extract Volume: 5000	(uL) Date Analyzed: 05/17/02
Injection Volume: 2.0(uL)	Dilution Factor: 1.0
C C Cleanup: (Y/N) N pH:	Sulfur Cleanup: (Y/N) N
CAS NO. COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L Q
12674-11-2Aroclor-1016_ 11104-28-2Aroclor-1221_ 11141-16-5Aroclor-1232_ 53469-21-9Aroclor-1242_ 12672-29-6Aroclor-1248_ 11097-69-1Aroclor-1254_ 11096-82-5Aroclor-1260_	0.50 U 1.0 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U

MINON

June 20, 2002 Compliance Sample Laboratory Results

SW-846

1-CC

CLASSICAL CHEMISTRY ANALYSES DATA SHEET

EPA SAMPLE NO.

l b Name: (CompuChem			Contract:				EFF	LUENT
_	LIBRTY		Case No.:		NRAS No.:				
G No.:	RG1024	_				_			
Matrix (soi	l/water):	WATER			La	b Sam	ple II	D: RG1024-	1
ite Receiv	ed: 6/21/02				*	Solid	s: <u>0</u>	.00	
	Concent	ration Unit	ts (mg/	L or mg/kg dry	weight	:) :	mg/	/L	
	PARAMETE	₹		CONCENTRATION	С	Ω	м	DATE ANALYZED]
	TSS]	1.20	1			6/25/02	- -
	рН			7.13	1			6/28/02	71

9/1/20/00

Comments:	
	2

FINAL REPORT OF ANALYSES

COMPUCHEM

Attn: DIANE BYRD 501 MADISON AVENUE CARY, NC 27513-

REPORT DATE: 06/28/02

SAMPLE NUMBER- 198141 SAMPLE ID- EFFLUENT

DATE SAMPLED- 06/20/02

DATE RECEIVED- 06/21/02 SAMPLER- NOT SPECIFIED

TIME RECEIVED- 1418

DELIVERED BY- CHRIS BRAND

Page 1 of 1

PROJECT NAME : ACS-89

ANALYSIS

ANALYSIS

METHOD DATE

BY RESULT UNITS

PQL

SAMPLE MATRIX- WW

RECEIVED BY- ALT

TIME SAMPLED- 1400

BIOCHEMICAL OXYGEN DEMAND

EPA 405.1 06/21/02 RCB

24 mg/L

2

QL = Practical Quantitation Limit
Results followed by the letter J are estimated concentrations.

NC DENR CERTIFICATIONS: DWQ - 96; PUBLIC WATER SUPPLY - 37724

LABORATORY DIRECTOR

1/30/00

SW846 METALS

INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

EFFLUENT

ab Name:	COMPUCHE	GM	Cont	ract:		<u> </u>		
ab Code:	LIBRTY	_ Case N	o.: s	AS No.:		SDG 1	No.:	RG1024
trix (so	il/water)	: WATER		Lab Sample	ID:	RG102	24-1	
evel (low	/med) ·	LOW		Date Receiv				
3VET (10"	, mea, .	<u> </u>		Dace Veceta	eu.	0/21/	-02	
Solids:	0.0	_						
				·-			110/	
		Concent	ration Units (ug.	/L or mg/kg dry	W61	gnt):	06/	<u> </u>
		CAS No.	Analyte	Concentration	C	Ω	М]
		7429-90-5	Aluminum	122		 	P	INB
		7440-38-2	Arsenic	4.2	Ū	i	P	
		7440-39-3		93.6		i	P	<u> </u>
		7440-36-0	Antimony	1.6		i	P	<u>.</u>
		7440-41-7	Beryllium	0.20	 -	<u> </u>	P	j
		7440-43-9	Cadmium	0.30	บ	1	P	-
		7440-70-2	Calcium	78400	T	<u> </u>	P	ĺ
		7440-47-3	Chromium	0.50	U	1	P	Ī
		7440-48-4	Cobalt	1.4	В		P	IUB
		7440-50-8	Copper	4.1	B	1	P	IUB
		7439-89-6	Iron	11.4	U	1	P	Ī
		7439-92-1	Lead	1.3	U	1	P	Ì
		7439-95-4	Magnesium	30600		1	P	ĺ
		7439-96-5	Manganese	24.5]	P	İ
		7439-97-6	Mercury	0.64	טן	1	CV	
		7440-02-0	Nickel	31.8		1	P	İ
		7440-09-7	Potassium	16600	1		P	
		7782-49-2	Selenium	2.2	ט		P	
		7440-22-4	Silver	0.70	ע		P	
		7440-23-5	Sodium	360000	<u> </u>	<u> </u>	P	l
		7440-28-0	Thallium	3.3		<u>l</u>	P	
		7440-62-2	Vanadium	2.0		<u> </u>	P	IUB
		7440-66-6	Zinc	3.9	B	<u> </u>	P	
								1/20/00
Color Bei	fore: COL	ORLESS	Clarity Before:	CLEAR	_ T	exture		
Color Aft	ter: COL	ORLESS	Clarity After:	CLEAR	_ A:	rtifac	ts: .	
Comments:	:							
								8 -

FORM 1 VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EFFLUENT

Lab Name: COMPUCHEM Method: 8260B

Lab Code: LIBRTY Case No.: SAS No.: SDG No.: RG1024

Matrix: (soil/water) WATER Lab Sample ID: RG1024-1

Sample wt/vol: 25 (g/ml) ML Lab File ID: RG1024-1A61

Level: (low/med) LOW Date Received: 06/21/02

% Moisture: not dec. _____ Date Analyzed: 06/29/02

GC Column: RTX-VMS ID: 0.18 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____(uL) Soil Aliquot Volume: _____(

CONCENTRATION UNITS:

CAS NO. COMPOUND (uq/L or uq/Kq) UG/L Q

CAB NO.	COMPOUND (dg/H of dg	/ kg/ 00/ L	¥
74-87-3	Chloromethane	0.2	J J
	Vinyl Chloride	0.5	ii)
74-83-9	Bromomethane	0.5	
	Chloroethane	0.5	
	1,1-Dichloroethene	0.5	
75-15-0	Carbon disulfide	0.5	II s
67-64-1	Acetone	4	BUBY
	Methylene Chloride	- i	BUBJ
156-60-5	trans-1,2-Dichloroethene	0.5	
75-34-3	1,1-Dichloroethane	0.5	
156-59-2	cis-1,2-Dichloroethene	0.5	U .
78-93-3	2-butanone	3	บันป
67-66-3	Chloroform	0.5	U
71-55-6	1,1,1-Trichloroethane	0.5	.บ
56-23-5	Carbon Tetrachloride	0.5	ប
	Benzene	0.5	บ
107-06-2	1,2-Dichloroethane	0.5	ַ ע
79-01-6	Trichloroethene	0.5	
78-87-5	1,2-Dichloropropane	0.5	
75-27-4	Bromodichloromethane	0.5	ַ U
10061-01-5-	cis-1,3-Dichloropropene	0.5	
	4-Methyl-2-pentanone	_{	U ,, a r
	Toluene	0.2	
10061-02-6-	trans-1,3-Dichloropropene	0.5	U ~
79-00-5	1,1,2-Trichloroethane	0.06	
127-18-4	Tetrachloroethene	0.5	
	2-hexanone	6	
124-48-1	Dibromochloromethane	0.5	1 1 1 LAT
108-90-7	Chlorobenzene	_} 0.09	JB UBJ
100-41-4	Ethylbenzene	0.5	
108-38-3	m,p-Xylene	_ 1	ן ט
	O-Xylene	0.5	
100-42-5	Styrene	0.5	ا ت ا
l	FORM T VOX	_	ll .

FORM I VOA

1/30/2

FORM 1 VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM	Method: 8260B	EFFLUENT
Lab Code: LIBRTY Case No.:	SAS No.: SDG	No.: RG1024
Matrix: (soil/water) WATER	Lab Sample ID:	RG1024-1
Sample wt/vol: 25 (g/ml) ML	Lab File ID:	RG1024-1A61
Level: (low/med) LOW	Date Received:	06/21/02
% Moisture: not dec.	Date Analyzed:	06/29/02
GC Column: RTX-VMS ID: 0.18 (mm)	Dilution Facto	or: 1.0
Soil Extract Volume:(uL)	Soil Aliquot V	Volume:(u
CAS NO. COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/I	
75-25-2Bromoform 79-34-51,1,2,2-Tetrack 541-73-11,3-Dichlorober 106-46-71,4-Dichlorober 95-50-11,2-Dichlorober 120-82-11,2,4-Trichloro 540-59-01,2-Dichloroet 1330-20-7Xylene (total)	nzenenzeneobenzenehene (total)	0.5 U 0.5 U 0.5 U 0.5 U 0.5 U 0.5 U

1/20/5

FORM I VOA

FORM 1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM	Method:	8270C	EFFLOEMI
Lab Code: LIBRTY	Case No.: SAS No	.: SDG	No.: RG1024
Matrix: (soil/water)	WATER	Lab Sample ID	: RG1024-1
Sample wt/vol:	1000 (g/mL) ML	Lab File ID:	RG1024-1B64
Level: (low/med)	LOW	Date Received	: 06/21/02
Moisture:	decanted: (Y/N)	Date Extracted	1:06/24/02
Concentrated Extract	Volume: 1000(uL)	Date Analyzed	: 06/25/02
Injection Volume:	1.0(uL)	Dilution Facto	or: 1.0
GPC Cleanup: (Y/N)	N pH:		
CAS NO.		ENTRATION UNITS: L or ug/Kg) UG/I	
106-44-5	Bis(2-chloroethyl)ethe 4-Methylphenol Isophorone bis(2-ethylhexyl)Phtha		9.6 U 10 U 10 U 4 JB

1/20/00

FORM I SV

8270C

FORM 1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Na	ame: COMPUCHEM		Method: 82	70C	EFFLUENT
Lab Co	ode: LIBRTY	Case No.:	SAS No.:	SDG	No.: RG1024
Matrix	k: (soil/water)	WATER	L	ab Sample ID:	RG1024-1
Sample	e wt/vol:	1000 (g/mL) ML	L	ab File ID:	RG1024-1A64
Level:	(low/med)	LOW	Da	ate Received:	06/21/02
% Mois	sture:	decanted: (Y/N)_	Da	ate Extracted	:06/25/02
Concer	ntrated Extract	Volume: 1000(uL) Da	ate Analyzed:	06/26/02
Inject	ion Volume:	1.0(uL)	Di	ilution Facto	r: 1.0
GPC Cl	eanup: (Y/N)	N pH:			
	CAS NO.	COMPOUND		RATION UNITS: r ug/Kg) UG/L	Q
	87-86-5	Pentachlorophe	nol		1 U

1/20/02

EFFLUENT Lab Name: COMPUCHEM Contract: 8082 Lab Code: LIBRTY Case No.: SAS No.: SDG No.: RG1024 Matrix: (soil/water) WATER Lab Sample ID: RG1024-1 Sample wt/vol: 1000 (g/mL) ML Lab File ID: decanted: (Y/N)___ % Moisture: Date Received: 06/21/02 Extraction: (SepF/Cont/Sonc) SEPF Date Extracted:06/26/02 Concentrated Extract Volume: 5000(uL) Date Analyzed: 06/27/02 Injection Volume: 2.0(uL) Dilution Factor: 1.0 Sulfur Cleanup: (Y/N) N GPC Cleanup: (Y/N) N pH: CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/L 0 0.50 U 12674-11-2----Aroclor-1016 11104-28-2----Aroclor-1221 1.0 0 11141-16-5-----Aroclor-1232 0.50 U uJ 0.50 U 53469-21-9-----Aroclor-1242 0.50 12672-29-6-----Aroclor-1248 U 11097-69-1-----Aroclor-1254 0.50 U 11096-82-5----Aroclor-1260 0.50 U

Washing.

APPENDIX B HISTORICAL WETLAND SEDIMENT ANALYTICAL DATA (May 1996)

TABLE 8 PCB ORGANICS ANALYSIS SUMMARY

Soil/Sediment Samples WETLAND INVESTIGATION

AMERICAN CHEMICAL SERVICES, INC.
GRIFFITH, INDIANA

SAMPLE ID		Detected PCBs				
	aroclor-1248	aroclor-1254	aroclor-1260			
SD17	58 JP	150 P	140	348		
SD18	••	••				
SD18-91		••	• •			
SD19	13 JP	36 JP	16 JP	65		
SD20		79 J	180 P	259		
SD21	1,300 JP	8,700	3,100 P	13,100		
SD22	560 JP	3,600	1,700	5,860		
SD22-91	270 JP	1,800	830	2,900		
SD23	770 JP	4,000	1,900	6,670		
SD24						
SD25	••	46 J		46		
SD26	320 JP	1,700	1,900	3,920		
SD27	48 J	190 P	270 P	508		
SD28	220 J	1,200	970 P	2,390		
SD29	180	380 P	330 P	890		
SD29-91	84 P	450 P	570 P	1,104		
SD30	74 P	570 P	390	1,034		
SD31	61 JP	600	240 P	901		
SD32	35 JP	79 P	73	187		
SD33	27,000 P	63,000 P	35,000 P	125,000		
SD34		14 JP	13 JP	27		
SD35	2,700 JP	8,100 P	6,200 P	17,000		
SD36		37 JP		37		
SD37						
SD38	30 JP	99 JP	100 P	229		

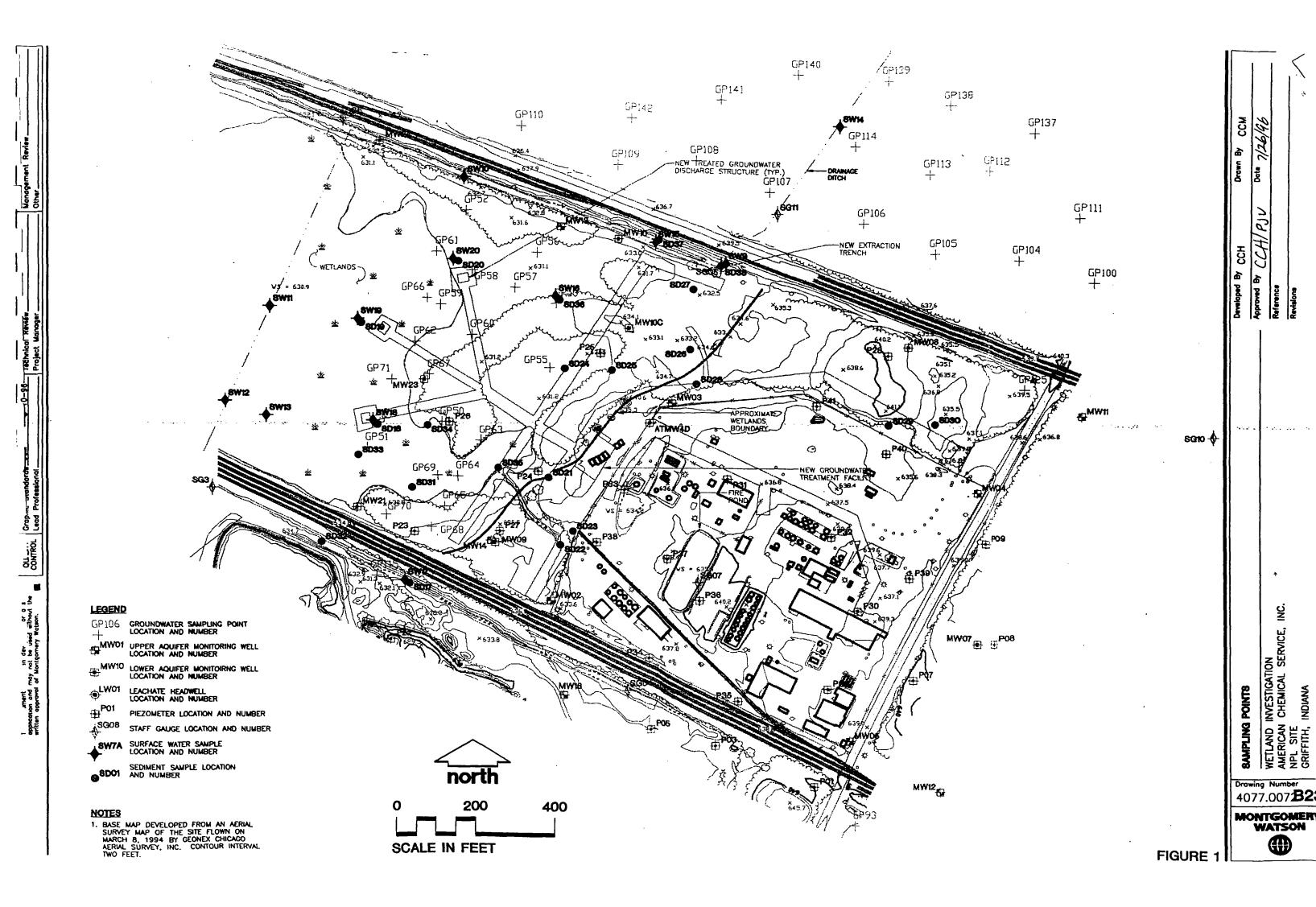
Notes:

- 1. All results expressed in micrograms per kilogram (ug/kg).
- 2. "--" = compound was not detected above the quantitation limit
- 3. "J" = indicates an estimated concentration between the quantitation limit and the method detection limit
- 4. "P" = This flag is used for pesticide/arochlor target analyte when there is a greater than 25 percent difference for detected concentrations between the two GC columns.

CCH/cch/SCI

j:\4077\0090\wetland\lab-data.xlw

PEST - soil



APPENDIX C SEDIMENT ANALYTICAL DATA

June 5, 2002 Sediment Sample Laboratory Results

GC EXTRACTABLE ORGANICS ANALYSIS DATA SHEET

COMPOUND

CAS NO.

EDA SAMPLE NO.

GWTP-003

0

Lab Name: COMPUCHEM Contract: 8082 Case No.: Lab Code: LIBRTY SAS No.: SDG No.: G2231 Matrix: (soil/water) SOIL Lab Sample ID: G2231-1 Lab File ID: Sample wt/vol: 30.0 (g/mL) G % Moisture: 37 decanted: (Y/N) Y Date Received: 06/06/02 Extraction: (SepF/Cont/Sonc) SONC Date Extracted:06/06/02 Concentrated Extract Volume: Date Analyzed: 06/10/02 5000 (uL) Injection Volume: 2.0(uL) Dilution Factor: 1.0 GPC Cleanup: Sulfur Cleanup: (Y/N) Y (Y/N) N pH:

7/20/00

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

GC EXTRACTABLE ORGANICS ANALYSIS DATA SHEET

EDA SAMPLE NO.

Lab	Name:	COMPUCHEM	Contract:	8082	GW1P-004

SDG No.: G2231 Lab Code: LIBRTY Case No.: SAS No.:

Matrix: (soil/water) SOIL Lab Sample ID: G2231-2

Lab File ID: Sample wt/vol: 30.0 (g/mL) G

Date Received: 06/06/02 % Moisture: 34 decanted: (Y/N) Y

Date Extracted:06/06/02 Extraction: (SepF/Cont/Sonc) SONC

Concentrated Extract Volume: 5000(uL) Date Analyzed: 06/10/02

Dilution Factor: 1.0 Injection Volume: 2.0(uL)

Sulfur Cleanup: (Y/N) Y GPC Cleanup: (Y/N) N pH:___

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG 0 49 U 12674-11-2-----Aroclor-1016 11104-28-2----Aroclor-1221 64 U 11141-16-5-----Aroclor-1232 49 U 34 U 53469-21-9-----Aroclor-1242 34 U 12672-29-6-----Aroclor-1248 34 11097-69-1-----Aroclor-1254 U U 11096-82-5-----Aroclor-1260 49

CONCENTRATION UNITS:

APPENDIX D CATALYTIC OXIDIZER OFF-GAS ANALYTICAL DATA

April 26, 2002 Off-Gas Sample (Round 1) Laboratory Results

AIR TOXICS LTD.

SAMPLE NAME: ACS-ME106-EF1-001A

ID#: 0204587A-01A

EPA METHOD TO-14 GC/MS FULL SCAN

130 12 mg	05000 2		Page + - eagle +	Turn Proper	
KII - Talkisee	<u></u>		Marque en afternación de P. C.		
Compound	Rpt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)	
Chloromethane	6.7	14	240	500	
Vinyl Chloride	6.7	17	350	900	
Bromomethane	6.7	26	Not Detected	Not Detected	
Chloroethane	6.7	18	170	460	
1,1-Dichloroethene	6.7	27	24	95	
Methylene Chloride	6.7	24	440	1500	
1,1-Dichloroethane	6.7	28	74	310	
cis-1,2-Dichloroethene	6.7	27	880	3600	
Chloroform	6.7	33	Not Detected	Not Detected	
1,1,1-Trichloroethane	6.7	37	13	72	
Carbon Tetrachloride	6.7	43	Not Detected	Not Detected	
Benzene	6.7	22	1500	4800	
1,2-Dichloroethane	6.7	28	32	130	
Trichloroethene	6.7	36	7.1	39	
1,2-Dichloropropane	6.7	31	7.4	35	
cis-1,3-Dichloropropene	6.7	31	Not Detected	Not Detected	
Toluene	6.7	26	750	2800	
trans-1,3-Dichloropropene	6.7	31	Not Detected	Not Detected	
1,1,2-Trichloroethane	6.7	37	2.0 J J	11 J	
Tetrachloroethene	6.7	46	5.3 J	37 J	
Chlorobenzene	6.7	31	110	540	
Ethyl Benzene	6.7	30	100	460	
m,p-Xylene	6.7	30	480	2100	
o-Xylene	6.7	30	160	730	
Styrene	6.7	29	19	84	
1,1,2,2-Tetrachloroethane	6.7	47	Not Detected	Not Detected	
Acetone	27	65	88	210	
Carbon Disulfide	27	85	Not Detected	Not Detected	
rans-1,2-Dichloroethene	27	110	54	220	
2-Butanone (Methyl Ethyl Ketone)	27	80	27	80	
Bromodichloromethane	27	180	Not Detected	Not Detected	
4-Methyl-2-pentanone	27	110	21 J	86 J	
2-Hexanone	27	110	Not Detected	Not Detected	
Dibromochloromethane	27	230	Not Detected	Not Detected	
Bromoform	27	280	Not Detected	Not Detected	

J = Estimated value.



AIR TOXICS LTD.

SAMPLE NAME: ACS-ME106-EF1-001A

ID#: 0204587A-01A

EPA METHOD TO-14 GC/MS FULL SCAN

set of an	9200		Posts of a difference of Weight	
office to peak			Andrews medium is reference	
			Method	
Surrogates	%Recovery		Limits	
1,2-Dichloroethane-d4		98	70-130	
Toluene-d8		96	70-130	
4-Bromofluorobenzene		92	70-130	

AIR TOXICS LTD.

SAMPLE NAME: ACS-ME106-IN1-001A

ID#: 0204587A-02A

EPA METHOD TO-14 GC/MS FULL SCAN

The North	th 1925 w		87.00 - 0 0.01/2.0	production of the production
	18.9		27.00. 00 00000	As Sec.
Compound	Rpt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)
Chloromethane	64	130	Not Detected	Not Detected
Vinyl Chloride	64	160	2900	7500
Bromomethane	64	250	Not Detected	Not Detected
Chloroethane	64	170	2200	5900
1,1-Dichloroethene	64	260	24 J 🏅	96 J
Methylene Chloride	64	220	3100	11000
1,1-Dichloroethane	64	260	1000	4300
cis-1,2-Dichloroethene	64	260	10000	41000
Chloroform	64	320	Not Detected	Not Detected
1,1,1-Trichloroethane	64	350	270	1500
Carbon Tetrachloride	64	410	Not Detected	Not Detected
Benzene	64	210	14000	45000
1.2-Dichloroethane	64	260	310	1300
Trichloroethene	64	350	65	360
1,2-Dichloropropane	64	300	110	510
cis-1,3-Dichloropropene	64	290	Not Detected	Not Detected
Toluene	64	240	10000	38000
trans-1,3-Dichloropropene	64	290	Not Detected	Not Detected
1,1,2-Trichloroethane	64	350	20 J 👅	110 J
Tetrachloroethene	64	440	23 J T	160 J
Chlorobenzene	64	300	920	4300
Ethyl Benzene	64	280	1600	7200
m,p-Xylene	64	280	7700	34000
o-Xylene	64	280	2500	11000
Styrene	64	270	Not Detected	Not Detected
1,1,2,2-Tetrachloroethane	64	440	Not Detected	Not Detected
Acetone	250	610	610	1500
Carbon Disulfide	250	800	Not Detected	Not Detected
trans-1,2-Dichloroethene	250	1000	Not Detected	Not Detected
2-Butanone (Methyl Ethyl Ketone)	250	760	340	1000
Bromodichloromethane	250	1700	Not Detected	Not Detected
4-Methyl-2-pentanone	250	1000	410	1700
2-Hexanone	250	1000	Not Detected	Not Detected
Dibromochloromethane	250	2200	Not Detected	Not Detected
Bromoform	250	2700	Not Detected	Not Detected

J = Estimated value.



SAMPLE NAME: ACS-ME106-IN1-001A

ID#: 0204587A-02A

EPA METHOD TO-14 GC/MS FULL SCAN

The end	1, 2977		(#151) en l'orghisteriffer (1524)
			Project to Majorage Commission
Surrogates		%Recovery	Method Limits
1,2-Dichloroethane-d4		98	70-130
Toluene-d8		96	70-130
4-Bromofluorobenzene		92	70-130

SAMPLE NAME: ACS-ME106-IN2-001A

ID#: 0204587A-03A

EPA METHOD TO-14 GC/MS FULL SCAN

170 (2 1 % 200) 18 (1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	6014 (1873) 380)		galaga i servici se dagar Negari sebili dagaraga	
Compound	Rpt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)
Chloromethane	340	710	Not Detected	Not Detected
Vinyl Chloride	340	880	7300	19000
Bromomethane	340	1300	Not Detected	Not Detected
Chloroethane	340	910	6600	18000
1,1-Dichloroethene	340	1400	410	1600
Methylene Chloride	340	1200	88000	310000
1,1-Dichloroethane	340	1400	20000	84000
cis-1,2-Dichloroethene	340	1400	9700	39000
Chloroform	340	1700	2000	10000
1,1,1-Trichloroethane	340	1900	53000	300000
Carbon Tetrachloride	340	2200	Not Detected	Not Detected ~
Benzene	340	1100	33000	110000
1,2-Dichloroethane	340	1400	1000	4300
Trichloroethene	340	1800	18000	100000
1,2-Dichloropropane	340	1600	270 J	1300 J
cis-1,3-Dichloropropene	340	1600	Not Detected	Not Detected
Toluene	340	1300	100000	400000
trans-1,3-Dichloropropene	340	1600	Not Detected	Not Detected
1,1,2-Trichloroethane	340	1900	Not Detected	Not Detected
Tetrachloroethene	340	2300	6800	47000
Chlorobenzene	340	1600	Not Detected	Not Detected
Ethyl Benzene	340	1500	7100	31000
m,p-Xylene	340	1500	26000	120000
o-Xylene	340	1500	7300	32000
Styrene	340	1500	Not Detected	Not Detected
1,1,2,2-Tetrachloroethane	340	2400	Not Detected	Not Detected
Acetone	1400	3300	16000	40000
Carbon Disulfide	1400	4300	Not Detected	Not Detected
trans-1,2-Dichloroethene	1400	5500	Not Detected	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1400	4100	15000	46000
Bromodichloromethane	1400	9300	Not Detected	Not Detected
4-Methyl-2-pentanone	1400	5700	3800	16000
2-Hexanone	1400	5700	Not Detected	Not Detected
Dibromochloromethane	1400	12000	Not Detected	Not Detected
Bromoform	1400	14000	Not Detected	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister



SAMPLE NAME: ACS-ME106-IN2-001A

ID#: 0204587A-03A

EPA METHOD TO-14 GC/MS FULL SCAN

At Weight			The set with region of briggs
			Burges Company to Child
Surrogates	_	%Recovery	Method Limits
1,2-Dichloroethane-d4		101	70-130
Toluene-d8		95	70-130
4-Bromofluorobenzene		94	70-130



SAMPLE NAME: ACS-ME106-EF1-001A

ID#: 0204587B-01A

EPA METHOD TO-13 GC/MS FULL SCAN

545 74 (o)	Artific to	• • • • • • • • • • • • • • • • • • •	Main in the appropriate the same
(c.)	6.4.1		Process of the Antonios of the
			Magnetina Chipere de Labor de Cater
	•	Rpt. Limit	Amount
Compound		(ug)	(ug)
Phenol		5.0	Not Detected
bis(2-Chloroethyl) Ether		1.0	Not Detected
2-Chlorophenol		5.0	Not Detected
1,3-Dichlorobenzene		1.0	Not Detected
1,4-Dichlorobenzene		1.0	1.0
1,2-Dichlorobenzene		1.0	1.6
2-Methylphenol (o-Cresol)		5.0	Not Detected
N-Nitroso-di-n-propylamine		1.0	Not Detected
4-Methylphenol		5.0	Not Detected
Hexachloroethane		1.0	Not Detected
Nitrobenzene		1.0	Not Detected
Isophorone		1.0	Not Detected
2-Nitrophenol		5.0	Not Detected
2,4-Dimethylphenol		5.0	Not Detected
bis(2-Chloroethoxy) Methane		1.0	Not Detected
2,4-Dichlorophenol		5.0	Not Detected
1,2,4-Trichlorobenzene		1.0	Not Detected
Naphthalene		1.0	0.97 J
4-Chloroaniline		10	Not Detected
Hexachlorobutadiene		1.0	Not Detected
4-Chloro-3-methylphenol		5.0	Not Detected
2-Methylnaphthalene		1.0	Not Detected
Hexachlorocyclopentadiene		20	Not Detected
2,4,6-Trichlorophenol		5.0	Not Detected
2,4,5-Trichlorophenol		5.0	Not Detected
2-Chloronaphthalene		1.0	Not Detected
2-Nitroaniline		10	Not Detected
Dimethylphthalate		5.0	Not Detected
Acenaphthylene		1.0	Not Detected
2,6-Dinitrotoluene		5.0	Not Detected
3-Nitroaniline		10	Not Detected
Acenaphthene		1.0	Not Detected
2,4-Dinitrophenol		20	Not Detected
4-Nitrophenol		20	Not Detected
2,4-Dinitrotoluene		5.0	Not Detected
Dibenzofuran		1.0	Not Detected
Diethylphthalate		5.0	Not Detected
Fluorene		1.0	Not Detected
4-Chlorophenyl-phenyl Ether		1.0	Not Detected
4-Nitroaniline		10	Not Detected
4,6-Dinitro-2-methylphenol		10	Not Detected
+,0-Dinitio-2-inethylphenol		.0	1101 Deleoted



SAMPLE NAME: ACS-ME106-EF1-001A

ID#: 0204587B-01A

EPA METHOD TO-13 GC/MS FULL SCAN

the state of the s	A Company of the Comp	
7 H. 21-00	a) 414	TREATE LAST SLEETES AND CARREST AND SECTION OF CONTROL
Section 1997		Principles Systems of the first
		disher to the designing of the better
		· · · · · · · · · · · · · · · · · · ·

	Rpt. Limit	Amount
Compound	(ug)	(ug)
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	1.2 J
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected •
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube: VOST

Surrogates	%Recovery	Method Limits
2-Fluorophenol	85	50-150
Phenol-d5	95	50-150
Nitrobenzene-d5	87	50-150
2-Fluorobiphenyl	90	60-120
2,4,6-Tribromophenol	84	50-150
Terphenyl-d14	106	60-120



SAMPLE NAME: ACS-ME106-IN1-001A

ID#: 0204587B-02A

EPA METHOD TO-13 GC/MS FULL SCAN

1 H 1 S D	with the		#Zac. of gettlessing acceptant
	,03		Supplied the converted process of \$10.000
			share ensettying comment of steels
		Rpt. Limit	Amount
Compound		(ug)	(ug)
Phenol		5.0	Not Detected
bis(2-Chloroethyl) Ether		1.0	Not Detected
2-Chlorophenol		5.0	Not Detected
1,3-Dichlorobenzene		1.0	0.88 J T
1,4-Dichlorobenzene		1.0	12
1,2-Dichlorobenzene		1.0	21
2-Methylphenol (o-Cresol)		5.0	Not Detected
N-Nitroso-di-n-propylamine		1.0	Not Detected
4-Methylphenol		5.0	Not Detected
Hexachloroethane		1.0	Not Detected
Nitrobenzene		1.0	Not Detected
Isophorone		1.0	Not Detected
2-Nitrophenol		5.0	Not Detected
2,4-Dimethylphenol		5.0	Not Detected
bis(2-Chloroethoxy) Methane		1.0	Not Detected
2,4-Dichlorophenol		5.0	Not Detected
1,2,4-Trichlorobenzene		1.0	0.75 J
Naphthalene		1.0	18
4-Chloroaniline		10	Not Detected
Hexachlorobutadiene		1.0	Not Detected
4-Chloro-3-methylphenol	······································	5.0	Not Detected
2-Methylnaphthalene		1.0	5.4
Hexachlorocyclopentadiene		20	Not Detected
2,4,6-Trichlorophenol		5.0	Not Detected
2,4,5-Trichlorophenol		5.0	Not Detected
2-Chloronaphthalene		1.0	Not Detected
2-Nitroaniline		10	Not Detected
Dimethylphthalate		5.0	Not Detected
Acenaphthylene		1.0	Not Detected
2,6-Dinitrotoluene		5.0	Not Detected
3-Nitroaniline		10	Not Detected
Acenaphthene		1.0	Not Detected
2,4-Dinitrophenol		20	Not Detected
4-Nitrophenol		20	Not Detected
2,4-Dinitrotoluene		5.0	Not Detected
Dibenzofuran		1.0	Not Detected
Diethylphthalate		5.0	Not Detected
Fluorene		1.0	Not Detected
4-Chlorophenyl-phenyl Ether		1.0	Not Detected
4-Nitroaniline		10	Not Detected
4,6-Dinitro-2-methylphenol		10	Not Detected
7,0-Dinitio-2-methylphenol			1101 00100100



SAMPLE NAME: ACS-ME106-IN1-001A

ID#: 0204587B-02A

EPA METHOD TO-13 GC/MS FULL SCAN

april 6 pers		Commence of the Commence of th
		The constant of the constant of the $\mathcal{M}_{\mathcal{A}}$
Compound	Rpt. Limit (ug)	Amount (ug)
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	2.0 J
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube: VOST

••		Method	
Surrogates	%Recovery	Limits	
2-Fluorophenol	66	50-150	
Phenol-d5	81	50-150	
Nitrobenzene-d5	83	50-150	
2-Fluorobiphenyl	89	60-120	
2,4,6-Tribromophenol	78	50-150	
Terphenyl-d14	105	60-120	



SAMPLE NAME: ACS-ME106-IN2-001A

ID#: 0204587B-03A

EPA METHOD TO-13 GC/MS FULL SCAN

of the American	Deal Section	रक्षात्व स्थानका[दुर्द्दीराज्य स्थानाप्र
	102	dans of Asglyss 35 m
		Digital on Englishing to a section
	Rpt. Limit	Amount
Compound	(ug)	(ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	Not Detected
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	Not Detected
1,4-Dichlorobenzene	1.0	Not Detected
1,2-Dichlorobenzene	1.0	3.9
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	Not Detected
Naphthalene	1.0	6.2
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	Not Detected
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
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SAMPLE NAME: ACS-ME106-IN2-001A

ID#: 0204587B-03A

EPA METHOD TO-13 GC/MS FULL SCAN

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NESS. TRANSPOR		$\tau_{i,i+1}$, we couplify $\mu(i,i+1)$, $\pi(i,i+1)$
		tion, or being exert, with the
	Rpt. Limit	Amount
Compound	(ug)	(ug)
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
ois(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
ndeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected
Container Type: XAD Tube: VOST		•• ••
Sandaa	9/ Pageyen:	Method Limits
Surrogates	%Recovery	
-Fluorophenol	88	50-150
Phenol-d5	99	50-150
Nitrobenzene-d5	94	50-150
2-Fluorobiphenyl	96	60-120
2,4,6-Tribromophenol	86	50-150
Ferphenyl-d14	113	60-120

May 22, 2002 Off-Gas Sample (Round 2) Laboratory Results

SAMPLE NAME: ACS-ME106-EF1-002A

ID#: 0205485-01A

EPA METHOD TO-14 GC/MS FULL SCAN

17.8 (R.)	(D. 7)		(18th) or configurate of the following	
			Sens of Employees of the s	
Compound	Røt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount Amount (ppbv) (uG/m3)	
Chloromethane	0.70	1.4	17 36	
Vinyl Chloride	0.70	1.8	57 150	
Bromomethane	0.70	2.7	Not Detected Not Detected	d
Chloroethane	0.70	1.9	18 49	
1,1-Dichloroethene	0.70	2.8	6.1 25	
Methylene Chloride	0.70	2.4	49 170	
1,1-Dichloroethane	0.70	2.8	14 60	
cis-1,2-Dichloroethene	0.70	2.8	180 750	
Chloroform	0.70	3.4	0.54J /J 2.7J	
1,1,1-Trichloroethane	0.70	3.8	3.1 17	
Carbon Tetrachloride	0.70	4.4	Not Detected Not Detected	d
Benzene	0.70	2.2	240 770	
1,2-Dichloroethane	0.70	2.8	5.4 22	
Trichloroethene	0.70	3.8	0.82 4.5	
1,2-Dichloropropane	0.70	3.3	1.4 6.8	
cis-1,3-Dichloropropene	0.70	3.2	Not Detected Not Detected	d
Toluene	0.70	2.7	120 460	
trans-1,3-Dichloropropene	0.70	3.2	Not Detected Not Detected	d
1,1,2-Trichloroethane	0.70	3.8	0.61 J / J 3.4 J	
Tetrachloroethene	0.70	4.8	0.30 J / T 2.1 J	
Chlorobenzene	0.70	3.2	23 110	•••••
Ethyl Benzene	0.70	3.1	20 86	
m,p-Xylene	0.70	3.1	74 320	
o-Xylene	0.70	3.1	26 120	
Styrene	0.70	3.0	3.8 16	
1,1,2,2-Tetrachloroethane	0.70	4.8	0.33 J / J 2.3 J	
Acetone	2.8	6.7	92 220	
Carbon Disulfide	2.8	8.8	0.31J /J 0.97J	
trans-1,2-Dichloroethene	2.8	11	8.5 34	
2-Butanone (Methyl Ethyl Ketone)	2.8	8.3	34 100	
Bromodichloromethane	2.8	19	Not Detected Not Detected	d
4-Methyl-2-pentanone	2.8	12	15 63	
2-Hexanone	2.8	12	0.46J /5 1.9J	
Dibromochloromethane	2.8	24	Not Detected Not Detected	d
Bromoform	2.8	29	Not Detected / Not Detected	

J = Estimated value.

LH 6/24/02

SAMPLE NAME: ACS-ME106-EF1-002A

ID#: 0205485-01A

EPA METHOD TO-14 GC/MS FULL SCAN

7 p \$600	(0,-12)		Figure 1996 september 2000
A Company	141	_	Region as integral and appeals of
Surrogates		%Recovery	Method Limits
1,2-Dichloroethane-d4		113	70-130
Toluene-d8		99	70-130
4-Bromofluorobenzene		95	70-130

SAMPLE NAME: ACS-ME106-IN1-002A

ID#: 0205485-02A

EPA METHOD TO-14 GC/MS FULL SCAN

The state	vitto int		Berlin, et allegeringen	
	Rpt. Limit	Rpt. Limit	Amount	Amount
Compound	(ppbv)	(uG/m3)	(ppbv)	(uG/m3)
Chloromethane	13	28	Not Detected	Not Detected
Vinyl Chloride	13	35	540	1400
Bromomethane	13	53	Not Detected	Not Detected
Chloroethane	13	36	280	760
1,1-Dichloroethene	13	54	4.9 J	20 J
Methylene Chloride	13	47	420	1500
1,1-Dichloroethane	13	55	210	880
cis-1,2-Dichloroethene	13	54	2300	9400
Chloroform	13	66	3.8 J /J	19 J
1,1,1-Trichloroethane	13	74	54	300
Carbon Tetrachloride	13	86	Not Detected	Not Detected
Benzene	13	44	2200	7100
1,2-Dichloroethane	13	55	53	220
Trichloroethene	13	73	6.3 J /J	34 J
1,2-Dichloropropane	13	63	22	100
cis-1,3-Dichloropropene	13	62	Not Detected	Not Detected
Toluene	13	51	1700	6600
trans-1,3-Dichloropropene	13	62	Not Detected	Not Detected
1,1,2-Trichloroethane	13	74	7.2J 15	40 J
Tetrachloroethene	13	92	Not Detected	Not Detected
Chlorobenzene	13	63	200	940
Ethyl Benzene	13	59	410	1800
m,p-Xylene	13	59	1700	7500
o-Xylene	13	59	580	2600
Styrene	13	58	Not Detected	Not Detected
1,1,2,2-Tetrachloroethane	13	94	3.5 J /5	25 J
Acetone	54	130	1100	2600
Carbon Disulfide	54	170	Not Detected	Not Detected
trans-1,2-Dichloroethene	54	220	Not Detected	Not Detected
2-Butanone (Methyl Ethyl Ketone)	54	160	630	1900
Bromodichloromethane	54	360	Not Detected	Not Detected
4-Methyl-2-pentanone	54	220	440	1800
2-Hexanone	54	220	13 J	53 J
Dibromochloromethane	54	460	Not Detected	Not Detected
Bromoform	54	560	Not Detected / Ū	Not Detected

J = Estimated value.

6/24/02

SAMPLE NAME: ACS-ME106-IN1-002A

ID#: 0205485-02A

EPA METHOD TO-14 GC/MS FULL SCAN

7.4	+17-241		type, echaditisan koatsszeti
			Police of Might on Physics
Surrogates		%Recovery	Method Limits
1,2-Dichloroethane-d4		112	70-130
Toluene-d8		98	70-130
4-Bromofluorobenzene		95	70-130

SAMPLE NAME: ACS-ME106-IN2-002A

ID#: 0205485-03A

EPA METHOD TO-14 GC/MS FULL SCAN

ert dagen. No. 100	Qs; (1)		North Control	
Compound	Rpt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)
Chloromethane	9.0	19	Not Detected	Not Detected
Vinyl Chloride	9.0	24	410	1100
Bromomethane	9.0	36	Not Detected	Not Detected
Chloroethane	9.0	24	210	580
1,1-Dichloroethene	9.0	36	3.7 J /J	15 J
Methylene Chloride	9.0	32	360	1200
1,1-Dichloroethane	9.0	37	170	720
cis-1,2-Dichloroethene	9.0	36	1900	7800
Chloroform	9.0	45	3.1 J	16 J
1,1,1-Trichloroethane	9.0	50	40	220
Carbon Tetrachloride	9.0	58	1.9 J /5	12 J
Benzene	9.0	29	1800	5700
1,2-Dichloroethane	9.0	37	46	190
Trichloroethene	9.0	49	5.4 J /J	30 J
1,2-Dichloropropane	9.0	42	20	92
cis-1,3-Dichloropropene	9.0	42	Not Detected	Not Detected
Toluene	9.0	35	1400	5500
trans-1,3-Dichloropropene	9.0	42	Not Detected	Not Detected
1,1,2-Trichloroethane	9.0	50	6.0J /T	33 J
Tetrachloroethene	9.0	62	Not Detected	Not Detected
Chlorobenzene	9.0	42	170	810
Ethyl Benzene	9.0	40	350	1500
m,p-Xylene	9.0	40	1400	6400
o-Xylene	9.0	40	520	2300
Styrene	9.0	39	Not Detected	Not Detected
1,1,2,2-Tetrachloroethane	9.0	63	2.7 J	19 J
Acetone	36	87	1100	2600
Carbon Disulfide	36	110	Not Detected	Not Detected
trans-1,2-Dichloroethene	36	140	Not Detected	Not Detected
2-Butanone (Methyl Ethyl Ketone)	· 36	110	630	1900
Bromodichloromethane	36	250	Not Detected	Not Detected
4-Methyl-2-pentanone	36	150	440	1800
2-Hexanone	36	150	12J /J	52 J
Dibromochloromethane	36	310	Not Detected	Not Detected
Bromoform	36	380	Not Detected / U	Not Detected

J = Estimated value.

LH 6/24/02

SAMPLE NAME: ACS-ME106-IN2-002A

ID#: 0205485-03A

EPA METHOD TO-14 GC/MS FULL SCAN

THE GOOD	(11456)		of signal long organization from a long to the con-
		Print of Meridans	
Surrogates		%Recovery	Method Limits
1,2-Dichloroethane-d4		111	70-130
Toluene-d8		100	70-130
4-Bromofluorobenzene		96	70-130

SAMPLE NAME: ACS-ME106-EF1-002A

ID#: 0205477-01A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

File Name: k060616 Date of Collection: 5/22/02
Dil. Factor: 1.00 Date of Analysis: 6/6/02
Date of Extraction: 5/23/02

	Rpt. Limit	Amount
Compound	(ug)	(ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	Not Detected
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	Not Detected
1,4-Dichlorobenzene	1.0	Not Detected
1,2-Dichlorobenzene	1.0	Not Detected
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	Not Detected
Naphthalene	1.0	Not Detected
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	Not Detected
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected

SAMPLE NAME: ACS-ME106-EF1-002A

ID#: 0205477-01A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

File Name:	k060616	Date of Collection: 5/22/02
Dil. Factor:	1.00	Date of Analysis: 6/6/02
		Date of Extraction: 5/23/02

	Rpt. Limit	Amount
Compound	(ug)	(ug)
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
ndeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

Container Type: XAD Tube: VOST

Container Type, 70 to Tube: 100		Method	
Surrogates	%Recovery	Limits	
2-Fluorophenol	108	50-150	
Phenol-d5	99	50-150	
Nitrobenzene-d5	88	50-150	
2-Fluorobiphenyl	89	60-120	
2,4,6-Tribromophenol	95	50-150	
Terphenyl-d14	90	60-120	

SAMPLE NAME: ACS-ME106-IN1-002A

ID#: 0205477-02A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

File Name:	k060617	Date of Collection: 5/22/02
Dil. Factor:	1.00	Date of Analysis: 6/6/02
	·	Date of Extraction: 5/23/02

	Rpt. Limit	Amount
Compound	(ug)	(ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	0.87J 万
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	0.50 J
1,4-Dichlorobenzene	1.0	6.3
1,2-Dichlorobenzene	1.0	13
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	1.6
Naphthalene	1.0	29
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	9.4
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected

SAMPLE NAME: ACS-ME106-IN1-002A

ID#: 0205477-02A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

File Name: k060617 Date of Collection: 5/22/02
Dil. Factor: 1.00 Date of Analysis: 6/6/02
Date of Extraction: 5/23/02

	Rpt. Limit	Amount
Compound	(ug)	(ug)
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	0.85 J <i>「</i> ブ
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube: VOST

••		Method Limits	
Surrogates	%Recovery		
2-Fluorophenol	106	50-150	
Phenol-d5	104	50-150	
Nitrobenzene-d5	107	50-150	
2-Fluorobiphenyl	107	60-120	
2,4,6-Tribromophenol	87	50-150	
Terphenyl-d14	109	60-120	

12 x 102

SAMPLE NAME: ACS-ME106-IN2-002A

ID#: 0205477-03A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

File Name:	k060618	Date of Collection: 5/22/02
Dil. Factor:	1.00	Date of Analysis: 6/6/02
		Date of Extraction: 5/23/02

	Rpt. Limit	Amount
Compound	(ug)	(ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1 0	0.70 J 13
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	Not Detected
1,4-Dichlorobenzene	1.0	4.4
1,2-Dichlorobenzene	1.0	9.2
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	1.2
Naphthalene	1.0	23
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	7.2
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected

SAMPLE NAME: ACS-ME106-IN2-002A

ID#: 0205477-03A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

File Name:	k060618	Date of Collection: 5/22/02
Dil. Factor:	1.00	Date of Analysis: 6/6/02
		Date of Extraction: 5/23/02

	Rpt. Limit	Amount
Compound	(ug)	(ug)
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	0.81 J / 3
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube: VOST

Surrogates	%Recovery	Method Limits	
Surroyates	/arecovery		
2-Fluorophenol	86	50-150	
Phenol-d5	84	50-150	
Nitrobenzene-d5	86	50-150	
2-Fluorobiphenyl	87	60-120	
2,4,6-Tribromophenol	76	50-150	
Terphenyl-d14	90	60-120	

June 21, 2002 Off-Gas Sample (Round 3) Laboratory Results

SAMPLE NAME: ACS-ME 106-EF1-003A

ID#: 0206434AR1-01A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

a at	412.74 (1.3)	gradica.		e nga sila sila danés sa terta. Wanasi silangdayan sa sag	
Compound	Rpt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)	
Chloromethane	29	60	170	370	
Vinyl Chloride	29	75	290	750	
Bromomethane	29	110	Not Detected	Not Detected	
Chloroethane	29	77	130	340	
1,1-Dichloroethene	29	120	26J <i>i</i> S	110 J	
Methylene Chloride	29	100	210	750	
1,1-Dichloroethane	29	120	54	220	
cis-1,2-Dichloroethene	29	120	750	3000	
Chloroform	29	140	Not Detected	Not Detected	
1,1,1-Trichloroethane	29	160	12 J (66 J	
Carbon Tetrachloride	29	180	Not Detected	Not Detected	
Benzene	29	94	1400	4700	
1.2-Dichloroethane	29	120	Not Detected	Not Detected	
Trichloroethene	29	160	د/ د9.0	49 J	
1,2-Dichloropropane	29	140	7.7 J 1	36 J	
cis-1,3-Dichloropropene	29	130	Not Detected	Not Detected	
Toluene	29	110	570	2200	
trans-1,3-Dichloropropene	29	130	Not Detected	Not Detected	
1,1,2-Trichloroethane	29	160	Not Detected	Not Detected	
Tetrachloroethene	29	200	5.3 J 👃	36 J	
Chlorobenzene	29	130	89	420	
Ethyl Benzene	29	130	84	370	
m,p-Xylene	29	130	360	1600	
o-Xylene	29	130	120	540	
Styrene	29	120	22 J 1	97 J	
1,1,2,2-Tetrachloroethane	29	200	Not Detected	Not Detected	
Acetone	120	280	200	490	
Carbon Disulfide	120	360	Not Detected	Not Detected	
trans-1,2-Dichloroethene	120	460	86 J / J	350 J	
2-Butanone (Methyl Ethyl Ketone)	120	340	8300	25000	
Bromodichloromethane	120	780	Not Detected	Not Detected	
4-Methyl-2-pentanone	120	480	22 J (S	93 J	
2-Hexanone	120	480	Not Detected	Not Detected	
Dibromochloromethane	120	1000	Not Detected	Not Detected	
Bromoform	120	1200	Not Detected	Not Detected	

J = Estimated value.



SAMPLE NAME: ACS-ME 106-EF1-003A

ID#: 0206434AR1-01A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

	£14 () 4 (
Surrogates	%Reco	Method overy Limits
1,2-Dichloroethane-d4	10	5 70-130
Toluene-d8	10	5 70-130
4-Bromofluorobenzene	10	1 70-130

SAMPLE NAME: ACS-ME 106-IN1-003A

ID#: 0206434AR1-02A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

			80 pt - 20 (1944)	that with the terminal	
			The second section of the section of the s		
Compound	Rpt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)	
Chloromethane	67	140	Not Detected	Not Detected	
Vinyl Chloride	67	170	2600	6800	
Bromomethane	67	260	Not Detected	Not Detected	
Chloroethane	67	180	1700	4600	
1,1-Dichloroethene	67	270	28J /J	110 J	
Methylene Chloride	67	240	1700	5900	
1,1-Dichloroethane	67	280	890	3600	
cis-1,2-Dichloroethene	67	270	9400 , ,	38000	
Chloroform	67	330	15J /S	75 J	
1,1,1-Trichloroethane	67	370	300	1700	
Carbon Tetrachloride	67	430	Not Detected	Not Detected	
Benzene	67	220	15000	48000	
1,2-Dichloroethane	67	280	Not Detected	Not Detected	
Trichloroethene	67	360	63 J /	340 J	
1,2-Dichloropropane	67	310	78	370	
cis-1,3-Dichloropropene	67	310	Not Detected	Not Detected	
Toluene	67	260	8500	33000	
trans-1,3-Dichloropropene	67	310	Not Detected	Not Detected	
1,1,2-Trichloroethane	67	370	Not Detected	Not Detected	
Tetrachloroethene	67	460	19J /J	130 J	
Chlorobenzene	67	310	740	3500	
Ethyl Benzene	67	300	1700	7600	
m,p-Xylene	67	300	8900	39000	
o-Xylene	67	300	2800	12000	
Styrene	67	290	Not Detected	Not Detected	
1,1,2,2-Tetrachloroethane	67	470	Not Detected	Not Detected	
Acetone	270	650	1200	2800	
Carbon Disulfide	270	850	Not Detected	Not Detected	
trans-1,2-Dichloroethene	270	1100	60 J / J	240 J	
2-Butanone (Methyl Ethyl Ketone)	270	800	610	1800	
Bromodichloromethane	270	1800	Not Detected	Not Detected	
4-Methyl-2-pentanone	270	1100	490	2000	
2-Hexanone	270	1100	Not Detected	Not Detected	
Dibromochloromethane	270	2300	Not Detected	Not Detected	
2.2.3.11001110110110110110	_ -				

J = Estimated value.

H 7/24/12

SAMPLE NAME: ACS-ME 106-IN1-003A

ID#: 0206434AR1-02A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

30 2 mg		Same and Migrae Room of the first	
April 1990 April 1990	2.2		Flore of Springer of the
Surrogates		%Recovery	Method Limits
1,2-Dichloroethane-d4		102	70-130
Toluene-d8		102	70-130
4-Bromofluorobenzene		99	70-130

At 7/210/02

SAMPLE NAME: ACS-ME 106-IN2-003A

ID#: 0206434AR1-03A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

No extension of the control of the c	green gan	eze (je		e de la companya de l	
Compound	Rpt. Limit _(ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)	
Chloromethane	70	150	Not Detected	Not Detected	
Vinyl Chloride	70	180	2600	6700	
Bromomethane	70	280	Not Detected	Not Detected	
Chloroethane	70	190	1700	4500	
1,1-Dichloroethene	70	280	23 J 🗘	94 J	
Methylene Chloride	70	250	1600	5700	
1,1-Dichloroethane	70	290	860	3500	
cis-1,2-Dichloroethene	70	280	9200	37000	
Chloroform	70	350	Not Detected	Not Detected	
1,1,1-Trichloroethane	70	390	290	1600	
Carbon Tetrachloride	70	450	Not Detected	Not Detected	
Benzene	70	230	15000	48000	
1.2-Dichloroethane	70	290	Not Detected	Not Detected	
Trichloroethene	70	380	60 J / J	33 0 J	
1,2-Dichloropropane	70	330	78	370	
cis-1,3-Dichloropropene	70	320	Not Detected	Not Detected	
Toluene	70	270	8400	32000	
trans-1,3-Dichloropropene	70	320	Not Detected	Not Detected	
1,1,2-Trichloroethane	70	390	Not Detected	Not Detected	
Tetrachloroethene	70	490	17 J /	110 J	
Chlorobenzene	70	330	750	3500	
Ethyl Benzene	70	310	1700	7500	
m,p-Xylene	70	310	8700	38000	
o-Xylene	70	310	2700	12000	
Styrene	70	300	Not Detected	Not Detected	
1,1,2,2-Tetrachloroethane	70	490	Not Detected	Not Detected	
Acetone	280	680	1100	2700	
Carbon Disulfide	280	890	Not Detected	Not Detected	
trans-1,2-Dichloroethene	280	1100	Not Detected	Not Detected	
2-Butanone (Methyl Ethyl Ketone)	280	840	580	1800	
Bromodichloromethane	280	1900	Not Detected	Not Detected	
4-Methyl-2-pentanone	280	1200	420	1800	
2-Hexanone	280	1200	Not Detected	Not Detected	
Dibromochloromethane	280	2400	Not Detected	Not Detected	
Bromoform	280	3000	Not Detected	Not Detected	

J = Estimated value.

Container Type: 6 Liter Summa Canister



SAMPLE NAME: ACS-ME 106-IN2-003A

ID#: 0206434AR1-03A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

100 to q ²	All though	The second second	Romand Contiference of the sign Roman year endposition of the	
Surrogates		%Recovery	Method Limits	
1,2-Dichloroethane-d4		100	70-130	
Toluene-d8		100	70-130	
4-Bromofluorobenzene		97	70-130	

SAMPLE NAME: ACS-ME 106-EF1-003A

ID#: 0206434BR1-01A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

Mr. N. and	16 15 114		Date of April About the State	
			Rieks + Schrepsbarz + 1000	
			Company of Degraphs of the section	
		Rpt. Limit	Amount	
Compound		(ug)	(ug)	
Phenol		5.0	Not Detected	
bis(2-Chloroethyl) Ether		1.0	Not Detected	
2-Chlorophenol		5.0	Not Detected	
1,3-Dichlorobenzene		1.0	Not Detected	
1,4-Dichlorobenzene		1.0	0.75 J	
1,2-Dichlorobenzene		1.0	1.2	
2-Methylphenol (o-Cresol)		5.0	Not Detected	
N-Nitroso-di-n-propylamine		1.0	Not Detected	
1-Methylphenol		5.0	Not Detected	
Hexachloroethane		1.0	Not Detected	
Nitrobenzene		1.0	Not Detected	
sophorone		1.0	Not Detected	
2-Nitrophenol		5.0	Not Detected	
2,4-Dimethylphenol		5.0	Not Detected	
ois(2-Chloroethoxy) Methane		1.0	Not Detected	
2,4-Dichlorophenol		5.0	Not Detected	
,2,4-Trichlorobenzene		1.0	Not Detected	
Naphthalene		1.0	Not Detected	
l-Chloroaniline		10	Not Detected	
Hexachlorobutadiene		1.0	Not Detected	
I-Chloro-3-methylphenol	***************************************	5.0	Not Detected	
2-Methylnaphthalene		1.0	Not Detected	
Hexachlorocyclopentadiene		20	Not Detected	
2,4,6-Trichlorophenol		5.0	Not Detected	
2,4,5-Trichlorophenol		5.0	Not Detected	
R-Chloronaphthalene		1.0	Not Detected	
2-Nitroaniline		10	Not Detected	
Dimethylphthalate		5.0	Not Detected	
Acenaphthylene		1.0	Not Detected	
2,6-Dinitrotoluene		5.0	Not Detected	
		10	Not Detected	
3-Nitroaniline		1.0	Not Detected	
Acenaphthene		20	Not Detected	
2,4-Dinitrophenol		20	Not Detected	
4-Nitrophenol		5.0	Not Detected	
2,4-Dinitrotoluene				
Dibenzofuran		1.0	Not Detected 0.34 J	
Diethylphthalate		5.0		
luorene		1.0	Not Detected	
l-Chlorophenyl-phenyl Ether		1.0	Not Detected	
I-Nitroaniline		10 10	Not Detected	

LH 7/2/02

SAMPLE NAME: ACS-ME 106-EF1-003A

ID#: 0206434BR1-01A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

	x(x,t) > C(x,t)		क्षिमा का स्वादा में किया है।
	.1()		Apply see morning puts
			Apolic of Contraction of Especia
		Rpt. Limit	Amount
Compound		(ug)	(ug)
N-Nitrosodiphenylamine		10	Not Detected
4-Bromophenyl-phenyl Ether		1.0	Not Detected
Hexachlorobenzene		1.0	Not Detected
Pentachlorophenol		20	Not Detected
Phenanthrene		1.0	Not Detected
Anthracene		1.0	Not Detected
di-n-Butylphthalate		5.0	Not Detected
Fluoranthene		1.0	Not Detected
Pyrene		1.0	Not Detected
Butylbenzylphthalate		5.0	Not Detected
3,3'-Dichlorobenzidine		20	Not Detected
Chrysene		1.0	Not Detected
Benzo(a)anthracene		1.0	Not Detected
bis(2-Ethylhexyl)phthalate		5.0	6.1
Di-n-Octylphthalate		5.0	Not Detected
Benzo(b)fluoranthene		1.0	Not Detected
Benzo(k)fluoranthene		1.0	Not Detected
Benzo(a)pyrene		1.0	Not Detected
ndeno(1,2,3-c,d)pyrene		1.0	Not Detected
Dibenz(a,h)anthracene		1.0	Not Detected
Benzo(g,h,i)perylene		1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube: VOST

-		Method	
Surrogates	%Recovery	Limits	
2-Fluorophenol	105	50-150	
Phenol-d5	96	50-150	
Nitrobenzene-d5	90	50-150	
2-Fluorobiphenyl	91	60-120	
2,4,6-Tribromophenol	106	50-150	
Terphenyl-d14	94	60-120	

SAMPLE NAME: ACS-ME 106-IN1-003A

ID#: 0206434BR1-02A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

(fig. 6) 360 (1.0) feet		bonds of all places of the gro
		Barrier of Training Community
		Charles of the output services of the edition of
	Rpt. Limit	Amount
Compound	(ug)	(ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	Not Detected
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	Not Detected
1,4-Dichlorobenzene	1.0	3.4
1,2-Dichlorobenzene	1.0	6.1
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
sophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
is(2-Chloroethoxy) Methane	1.0	Not Detected
,4-Dichlorophenol	5.0	Not Detected
,2,4-Trichlorobenzene	1.0	0.66 J 🔏
Naphthalene	1.0	2.8
-Chloroaniline	10	Not Detected
	1.0	Not Detected
l-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	1.4
lexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
?-Chloronaphthalene	1.0	Not Detected
-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
cenaphthylene	1.0	Not Detected
:,6-Dinitrotoluene	5.0	Not Detected
-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
-Nitrophenol	20	Not Detected
,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	0.37 J
luorene	1.0	Not Detected
-Chlorophenyl-phenyl Ether	1.0	Not Detected
Chiorophenyi-phenyi Ether Nitroaniline	10	Not Detected
,-Nttroamme ,6-Dinitro-2-methylphenol	10	Not Detected

SAMPLE NAME: ACS-ME 106-IN1-003A

ID#: 0206434BR1-02A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

	1.47 to 6		ामक क्षेत्रका ाना (हिन्दिन्नि	
	1.1		there is greater than the state of	
			The state of the fig. following the asset of	
		Rpt. Limit	Amount	
Compound		(ug)	(ug)	
N-Nitrosodiphenylamine		10	Not Detected	
4-Bromophenyl-phenyl Ether		1.0	Not Detected	
Hexachlorobenzene		1.0	Not Detected	
Pentachlorophenol		20	Not Detected	
Phenanthrene		1.0	Not Detected	
Anthracene		1.0	Not Detected	••••
di-n-Butylphthalate		5.0	Not Detected	
Fluoranthene		1.0	Not Detected	
Pyrene		1.0	Not Detected	
Butylbenzylphthalate		5.0	Not Detected	
3,3'-Dichlorobenzidine		20	Not Detected	
Chrysene		1.0	Not Detected	
Benzo(a)anthracene		1.0	Not Detected	سيسو
bis(2-Ethylhexyl)phthalate		5.0	2.0 J	13
Di-n-Octylphthalate		5.0	Not Detected	
Benzo(b)fluoranthene		1.0	Not Detected	••••
Benzo(k)fluoranthene		1.0	Not Detected	
Benzo(a)pyrene		1.0	Not Detected	
Indeno(1,2,3-c,d)pyrene		1.0	Not Detected	
Dibenz(a,h)anthracene		1.0	Not Detected	
Benzo(g,h,i)perylene		1.0	Not Detected	
J = Estimated value.				
Container Type: XAD Tube: VOST				
			Method	
Surrogates		%Recovery	Limits	_
2-Fluorophenol		88	50-150	
Phenol-d5		91	50-150	
Nitrobenzene-d5		86	50-150	
2-Fluorobiphenyl		87	60-120	

LH 7/26/02

50-150

60-120

101

92

2,4,6-Tribromophenol

Terphenyl-d14

SAMPLE NAME: ACS-ME 106-IN2-003A

ID#: 0206434BR1-03A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

the Major Comment		May be seen to the section of the se
(f) (s)(6)		A part of a Metric of the Control of
		in part \$1000 ken as open in the later.
_	Rpt. Limit	Amount
Compound	(ug)	(ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	1.2
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	3.0
1,4-Dichlorobenzene	1.0	32
1,2-Dichlorobenzene	1.0	56
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
pis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	7.1
Naphthalene	1.0	27
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	15
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
I-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	0.43 J
Fluorene	1.0	Not Detected
I-Chlorophenyl-phenyl Ether	1.0	Not Detected
I-Nitroaniline	10	Not Detected
l,6-Dinitro-2-methylphenol	10	Not Detected

SAMPLE NAME: ACS-ME 106-IN2-003A

ID#: 0206434BR1-03A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

77 At 128 source	existe in:		भिन्ना विकासितिक विभिन्न	
Trip to me	(ii)		THE MANGEMENT WHITE	
			कार्यन्त्री अस्ति स्ट्रांस्ट्रांस	
	<u></u>	Rpt. Limit	Amount	
Compound		(ug)	(ug)	
N-Nitrosodiphenylamine		10	Not Detected	_
4-Bromophenyl-phenyl Ether		1.0	Not Detected	
Hexachlorobenzene		1.0	Not Detected	
Pentachlorophenol		20	Not Detected	
Phenanthrene		1.0	Not Detected	
Anthracene		1.0	Not Detected	
di-n-Butylphthalate		5.0	1.4 J	1.2
Fluoranthene		1.0	Not Detected	
Pyrene		1.0	Not Detected	
Butylbenzylphthalate		5.0	Not Detected	
3,3'-Dichlorobenzidine		20	Not Detected	
Chrysene		1.0	Not Detected	
Benzo(a)anthracene		1.0	Not Detected	1
bis(2-Ethylhexyl)phthalate		5.0	1.9 J	15
Di-n-Octylphthalate		5.0	Not Detected	
Benzo(b)fluoranthene		1.0	Not Detected	
Benzo(k)fluoranthene		1.0	Not Detected	
Benzo(a)pyrene		1.0	Not Detected	
indeno(1,2,3-c,d)pyrene		1.0	Not Detected	
Dibenz(a,h)anthracene		1.0	Not Detected	
Benzo(g,h,i)perylene		1.0	Not Detected	
J = Estimated value.				
Container Type: XAD Tube: VOST				
			Method	
Surrogates		%Recovery	Limits	
2-Fluorophenol		54	50-150	
Phenol-d5		85	50-150	

88

91

98

93

1H 7/21/02

50-150

60-120

50-150

60-120

Nitrobenzene-d5 2-Fluorobiphenyl

Terphenyl-d14

2,4,6-Tribromophenol

June 28, 2002 Off-Gas Sample (Round 4) Laboratory Results

SAMPLE NAME: ACS-ME 106-EF1-004A

ID#: 0206594A-01A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

in the state of th	100 (CS)		Ortography	AND THE SEC
Compound	Rpt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)
Chloromethane	7.3	15	180	370
Vinyl Chloride	7.3	19	440	1100
Bromomethane	7.3	29	Not Detected	Not Detected
Chloroethane	7.3	20	170	450
1,1-Dichloroethene	7.3	29	32	130
Methylene Chloride	7.3	26	130	450
1,1-Dichloroethane	7.3	30	58	240
cis-1,2-Dichloroethene	7.3	29	810	3300
Chloroform	7.3	36	1.2J /	6.0 J
1,1,1-Trichloroethane	7.3	40	15	84
Carbon Tetrachloride	7.3	47	Not Detected	Not Detected
Benzene	7.3	24	1700	5600
1,2-Dichloroethane	7.3	30	Not Detected	Not Detected
Trichloroethene	7.3	40	8.4	46
1,2-Dichloropropane	7.3	34	4.5 J /	21 J
cis-1,3-Dichloropropene	7.3	34	Not Detected	Not Detected
Toluene	7.3	28	640	2500
trans-1,3-Dichloropropene	7.3	34	Not Detected	Not Detected
1,1,2-Trichloroethane	7.3	40	Not Detected	Not Detected
Tetrachloroethene	7.3	50	6.9 J J	47 J
Chlorobenzene	7.3	34	80	370
Ethyl Benzene	7.3	32	93	410
m,p-Xylene	7.3	32	460	2000
o-Xylene	7.3	32	130	570
Styrene	7.3	32	22	94
1,1,2,2-Tetrachloroethane	7.3	51	Not Detected	Not Detected
Acetone	29	70	48	120
Carbon Disulfide	29	92	7.2 J 🞵	23 J
trans-1,2-Dichloroethene	29	120	54	220
2-Butanone (Methyl Ethyl Ketone)	29	88	32	97
Bromodichloromethane	29	200	Not Detected	Not Detected
4-Methyl-2-pentanone	29	120	10 J J	42 J
2-Hexanone	29	120	Not Detected	Not Detected
Dibromochloromethane	29	250	Not Detected	Not Detected
Bromoform	29	310	Not Detected	Not Detected

J = Estimated value.

1/2/e/02

SAMPLE NAME: ACS-ME 106-EF1-004A

ID#: 0206594A-01A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

eliating		তাওঁমতিক জন উত্তর্গত নাথেকে এক উপন্যান
Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	115	70-130
Toluene-d8	95	70-130
4-Bromofluorobenzene	93	70-130

SAMPLE NAME: ACS-ME 106-IN1-004A

ID#: 0206594A-02A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

Aloka Eli-	A 4		Detection (April)	150 <u>01</u> 30 3.70102
Compound	Rpt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)
Chloromethane	67	140	Not Detected	Not Detected
Vinyl Chloride	67	170	3900	10000
Bromomethane	67	260	Not Detected	Not Detected
Chloroethane	67	180	2000	5300
1,1-Dichloroethene	67	270	25 J /5	100 J
Methylene Chloride	67	240	860	3000
1,1-Dichloroethane	67	280	780	3200
cis-1,2-Dichloroethene	67	270	7700	31000
Chloroform	67	330	12J 15	57 J
1,1,1-Trichloroethane	67	370	340	1900
Carbon Tetrachloride	67	430	Not Detected	Not Detected
Benzene	67	220	14000	46000
1,2-Dichloroethane	67	280	Not Detected	Not Detected
Trichloroethene	67	360	68	370
1,2-Dichloropropane	67	310	60 J (Z	280 J
cis-1,3-Dichloropropene	67	310	Not Detected	Not Detected
Toluene	67	260	7600	29000
trans-1,3-Dichloropropene	67	310	Not Detected	Not Detected
1,1,2-Trichloroethane	67	370	Not Detected	Not Detected
Tetrachloroethene	67	460	31 J /J	220 J
Chlorobenzene	67	310	570	2700
Ethyl Benzene	67	300	1600	6800
m,p-Xylene	67	300	8400	37000
o-Xylene	67	300	2200	9600
Styrene	67	290	63 J /\(\sum_{\subset}\)	270 J
1,1,2,2-Tetrachloroethane	67	470	Not Detected	Not Detected
Acetone	270	650	350	850
Carbon Disulfide	270	850	Not Detected	Not Detected
trans-1,2-Dichloroethene	270	1100	Not Detected	Not Detected
2-Butanone (Methyl Ethyl Ketone)	270	800	200 J /5	610 J
Bromodichloromethane	270	1800	Not Detected	Not Detected
4-Methyl-2-pentanone	270	1100	240 J /J	1000 J
2-Hexanone	270	1100	Not Detected	Not Detected
Dibromochloromethane	270	2300	Not Detected	Not Detected
Bromoform	270	2800	Not Detected	Not Detected

J = Estimated value.

1/21d02

SAMPLE NAME: ACS-ME 106-IN1-004A

ID#: 0206594A-02A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

Cillian Holes	25.	OF DOCK UPLIED AND OF	
Surrogates	%Recovery	Method Limits	
1,2-Dichloroethane-d4	114	70-130	
Toluene-d8	94	70-130	
4-Bromofluorobenzene	95	70-130	

SAMPLE NAME: ACS-ME 106-IN2-004A

ID#: 0206594A-03A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

File Seni Sign (Sec.	ही। इ.स.		सम्बद्धाः स्टब्स् वेम्बस्य स्टिब्स	
Compound	Røt. Limit (ppbv)	Rpt. Limit (uG/m3)	Amount (ppbv)	Amount (uG/m3)
Chloromethane	67	140	Not Detected	Not Detected
Vinyl Chloride	67	170	4600	12000
Bromomethane	67	260	Not Detected	Not Detected
Chloroethane	67	180	2300	6200
1,1-Dichloroethene	67	270	29 J 🗘	120 J
Methylene Chloride	67	240	1200	4400
1,1-Dichloroethane	67	280	970	4000
cis-1,2-Dichloroethene	67	270	10000	41000
Chloroform	67	330	13 J 🔼	65 J
1,1,1-Trichloroethane	67	370	410	2300
Carbon Tetrachloride	67	430	Not Detected	Not Detected
Benzene	67	220	17000	55000
1,2-Dichloroethane	67	280	Not Detected	Not Detected
Trichloroethene	67	360	89	490
1,2-Dichloropropane	67	310	87	410
cis-1,3-Dichloropropene	67	310	Not Detected	Not Detected
Toluene	67	260	12000	45000
trans-1,3-Dichloropropene	67	310	Not Detected	Not Detected
1,1,2-Trichloroethane	67	370	Not Detected	Not Detected
Tetrachloroethene	67	460	ن/ 31 ا	210 J
Chlorobenzene	67	310	820	3800
Ethyl Benzene	67	300	2300	10000
m,p-Xylene	67	300	12000	53000
o-Xylene	67	300	3300	15000
Styrene	67	290	Not Detected	Not Detected
1,1,2,2-Tetrachloroethane	67	470	Not Detected	Not Detected
Acetone	270	650	400	970
Carbon Disulfide	270	850	Not Detected	Not Detected
trans-1,2-Dichloroethene	270	1100	Not Detected	Not Detected
2-Butanone (Methyl Ethyl Ketone)	270	800	240 J /	720 J
Bromodichloromethane	270	1800	Not Detected	Not Detected
4-Methyl-2-pentanone	270	1100	330	1400
2-Hexanone	270	1100	Not Detected	Not Detected
Dibromochloromethane	270	2300	Not Detected	Not Detected
Bromoform	270	2800	Not Detected	Not Detected

J = Estimated value.

1/26/02

SAMPLE NAME: ACS-ME 106-IN2-004A

ID#: 0206594A-03A

MODIFIED EPA METHOD TO-14 GC/MS FULL SCAN

	The state of the s	可以中心的 1000 1000 1000 1000 1000 1000 1000 10	
Surrogates	%Recovery	Method Limits	
1,2-Dichloroethane-d4	116	70-130	
Toluene-d8	110	70-130	
4-Bromofluorobenzene	95	70-130	

SAMPLE NAME: ACS-ME 106-EF1-004A

ID#: 0206594B-01A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

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			The state of the s
<u> 1900 y la la participa de la compania del compania de la compania de la compania del compania de la compania del la compania del la compania de la compania de la compania de la compania de la compania de la compania de la compania de la compania de la compania del la compania del la compania de la compania del la</u>	Rpt	Limit	Amount
Compound	(1	ıg)	(ug)
Phenol	5	5.0	Not Detected
bis(2-Chloroethyl) Ether	1	.0	Not Detected
2-Chlorophenol	5	5.0	Not Detected
1,3-Dichlorobenzene	1	.0	Not Detected
1,4-Dichlorobenzene	1	.0	1.5
1,2-Dichlorobenzene	1	.0	2.3
2-Methylphenol (o-Cresol)	5	0.0	Not Detected
N-Nitroso-di-n-propylamine	1	.0	Not Detected
4-Methylphenol	5	5.0	Not Detected
Hexachloroethane	1	.0	Not Detected
Nitrobenzene	1	.0	Not Detected
sophorone	1	.0	Not Detected
2-Nitrophenol	5	.0	Not Detected
2,4-Dimethylphenol	5	.0	Not Detected
ois(2-Chloroethoxy) Methane	1	.0	Not Detected
2,4-Dichlorophenol	5	.0	Not Detected
1,2,4-Trichlorobenzene	1	.0	0.38J / _
Naphthalene	1	.0	0.41 J / 🗸
4-Chloroaniline	1	0	Not Detected
Hexachlorobutadiene	1	.0	Not Detected
1-Chloro-3-methylphenol	5	.0	Not Detected
2-Methylnaphthalene	1	.0	Not Detected
-lexachlorocyclopentadiene	2	20	Not Detected
2,4,6-Trichlorophenol	5	.0	Not Detected
2,4,5-Trichlorophenol		.0	Not Detected
2-Chloronaphthalene	1	.0	Not Detected
2-Nitroaniline	1	0	Not Detected
Dimethylphthalate	5	.0	Not Detected
Acenaphthylene	1	.0	Not Detected
2,6-Dinitrotoluene	5	.0	Not Detected
3-Nitroaniline		0	Not Detected
Acenaphthene		.0	Not Detected
2,4-Dinitrophenol		0	Not Detected
I-Nitrophenol		0	Not Detected
2,4-Dinitrotoluene		.0	Not Detected
Dibenzofuran		.0	Not Detected
Diethylphthalate		.0	0.47 J
Fluorene		.0	Not Detected
I-Chlorophenyl-phenyl Ether		.0	Not Detected
I-Nitroaniline		0	Not Detected
1,6-Dinitro-2-methylphenol		0	Not Detected

SAMPLE NAME: ACS-ME 106-EF1-004A

ID#: 0206594B-01A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

Distriction of the control of the co	i i de la caracteria de		
San San San San San San San San San San	Rpt. Limit	Amount	
Compound	(ug)	(ug)	
N-Nitrosodiphenylamine	10	Not Detected	
1-Bromophenyl-phenyl Ether	1.0	Not Detected	
Hexachlorobenzene	1.0	Not Detected	
Pentachlorophenol	20	Not Detected	
Phenanthrene	1.0	Not Detected	
Anthracene	1.0	Not Detected	
di-n-Butylphthalate	5.0	Not Detected	
Fluoranthene	1.0	Not Detected	
Pyrene	1.0	Not Detected	
Butylbenzylphthalate	5.0	Not Detected	
3,3'-Dichlorobenzidine	20	Not Detected	
Chrysene	1.0	Not Detected	
Benzo(a)anthracene	1.0	Not Detected	
pis(2-Ethylhexyl)phthalate	5.0	Not Detected	
Di-n-Octylphthalate	5.0	Not Detected	
Benzo(b)fluoranthene	1.0	Not Detected	
Benzo(k)fluoranthene	1.0	Not Detected	
Benzo(a)pyrene	1.0	Not Detected	
ndeno(1,2,3-c,d)pyrene	1.0	Not Detected	
Dibenz(a,h)anthracene	1.0	Not Detected	
Benzo(g,h,i)perylene	1.0	Not Detected	

J = Estimated value.

Container Type: XAD Tube: VOST

Surrogates	%Recovery	Method Limits
2-Fluorophenol	101	50-150
Phenol-d5	95	50-150
Nitrobenzene-d5	80	50-150
2-Fluorobiphenyl	86	60-120
2,4,6-Tribromophenol	112	50-150
Terphenyl-d14	93	60-120

SAMPLE NAME: ACS-ME 106-IN1 004A

ID#: 0206594B-02A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

1900	EDE THE THE PARTY OF THE PARTY	Unico (Sollegion) 182802
	8.56	The state of the s
		Applications to 12/102
designed in section of the control o	Rpt. Limit	Amount
Compound	(ug)	(ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	1.7
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.7
1,4-Dichlorobenzene	1.0	27
1,2-Dichlorobenzene	1.0	46
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	6.2
Naphthalene	1.0	13
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	6.9
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
	1.0	Not Detected
Acenaphthene	20	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	5.0	Not Detected
2,4-Dinitrotoluene	······································	Not Detected
Dibenzofuran Diethylahthalata	1.0 5.0	0.52 J
Diethylphthalate		0.020
Fluorene	1.0	Not Detected
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected

SAMPLE NAME: ACS-ME 106-IN1 004A

ID#: 0206594B-02A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

		্রেক্টিরিয়েদ্র দে প্রদেশ কর্মনার করে। ভারতির বিদ্যালয় বিশ্বন
Compound	Rpt. Limit (ug)	Amount (ug)
N-Nitrosodiphenylamine	10	Not Detected
I-BromophenyI-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
li-n-Butylphthalate	5.0	1.2 J
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
sis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
ndeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Q = Exceeds Quality Control limits.
Container Type: XAD Tube: VOST

Surrogates	%Recovery	Method Limits
2-Fluorophenol	44 Q	50-150
Phenol-d5	70	50-150
Nitrobenzene-d5	77	50-150
2-Fluorobiphenyl	76	60-120
2,4,6-Tribromophenol	61	50-150
Terphenyl-d14	81	60-120

1/26/02

SAMPLE NAME: ACS-ME 106-IN2 004A

ID#: 0206594B-03A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

Circumstantian (for		
		क इंद्रोहर्स्छ । इसस
	Rpt. Limit	Amount
Compound	(ug)	(ug)
Phenol	5.0	Not Detected
ois(2-Chloroethyl) Ether	1.0	Not Detected
2-Chlorophenol	5.0	Not Detected
,3-Dichlorobenzene	1.0	Not Detected
,4-Dichlorobenzene	1.0	Not Detected
,2-Dichlorobenzene	1.0	Not Detected
?-Methylphenol (o-Cresol)	5.0	Not Detected
I-Nitroso-di-n-propylamine	1.0	Not Detected
-Methylphenol	5.0	Not Detected
lexachioroethane	1.0	Not Detected
litrobenzene	1.0	Not Detected
sophorone	1.0	Not Detected
-Nitrophenol	5.0	Not Detected
,4-Dimethylphenol	5.0	Not Detected
is(2-Chloroethoxy) Methane	1.0	Not Detected
,4-Dichlorophenol	5.0	Not Detected
,2,4-Trichlorobenzene	1.0	Not Detected
iaphthalene	1.0	Not Detected
-Chloroaniline	10	Not Detected
lexachlorobutadiene	1.0	Not Detected
-Chloro-3-methylphenol	5.0	Not Detected
-Methylnaphthalene	1.0	Not Detected
lexachlorocyclopentadiene	20	Not Detected
,4,6-Trichlorophenol	5.0	Not Detected
,4,5-Trichlorophenol	5.0	Not Detected
-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
cenaphthylene	1.0	Not Detected
,6-Dinitrotoluene	5.0	Not Detected
-Nitroaniline	10	Not Detected
cenaphthene	1.0	Not Detected
,4-Dinitrophenol	20	Not Detected
-Nitrophenol	20	Not Detected
,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
fluorene	1.0	Not Detected
-Chlorophenyl-phenyl Ether	1.0	Not Detected
-Nitroaniline	10	Not Detected
,6-Dinitro-2-methylphenol	10	Not Detected

SAMPLE NAME: ACS-ME 106-IN2 004A

ID#: 0206594B-03A

MODIFIED EPA METHOD TO-13 GC/MS FULL SCAN

FIOLEN Outcon	Dat	Dand (collection (270,02) Dand Analysts 7000 Dand Analysts 7000	
	Rpt. Limit	Amount	
Compound	(ug)	(ug)	
N-Nitrosodiphenylamine	10	Not Detected	
4-Bromophenyl-phenyl Ether	1.0	Not Detected	
Hexachlorobenzene	1.0	Not Detected	
Pentachlorophenol	20	Not Detected	
Phenanthrene	1.0	Not Detected	
Anthracene	1.0	Not Detected	
di-n-Butylphthalate	5.0	Not Detected	
Fluoranthene	1.0	Not Detected	
Pyrene	1.0	Not Detected	
Butylbenzylphthalate	5.0	Not Detected	
3,3'-Dichlorobenzidine	20	Not Detected	
Chrysene	1.0	Not Detected	
Benzo(a)anthracene	1.0	Not Detected	
bis(2-Ethylhexyl)phthalate	5.0	Not Detected	
Di-n-Octylphthalate	5.0	Not Detected	
Benzo(b)fluoranthene	1.0	Not Detected	
Benzo(k)fluoranthene	1.0	Not Detected	
Benzo(a)pyrene	1.0	Not Detected	
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected	
Dibenz(a,h)anthracene	1.0	Not Detected	
Benzo(g,h,i)perylene	1.0	Not Detected	
Container Type: XAD Tube: VOST			
Surragatas	% Pagayan	Method Limits	
Surrogates	%Recovery		
2-Fluorophenol	99	50-150	
Phenol-d5	89	50-150	
Nitrobenzene-d5	83	50-150	
2-Fluorobiphenyl	84	60-120	
2,4,6-Tribromophenol	92	50-150	

60-120

Terphenyl-d14

88